

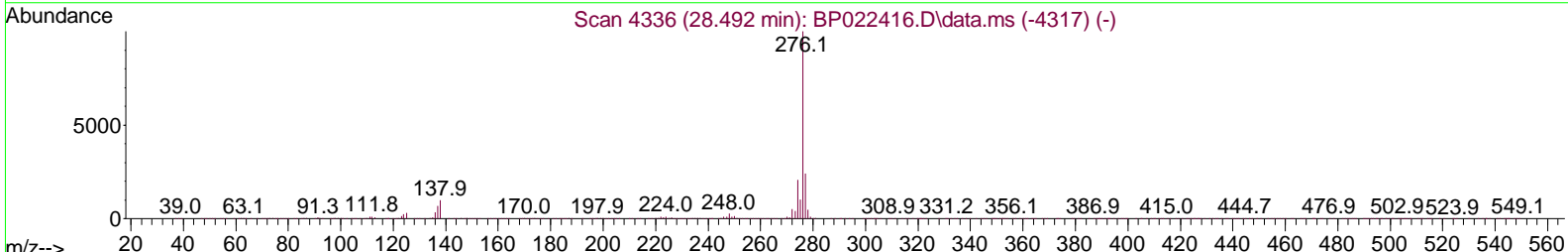
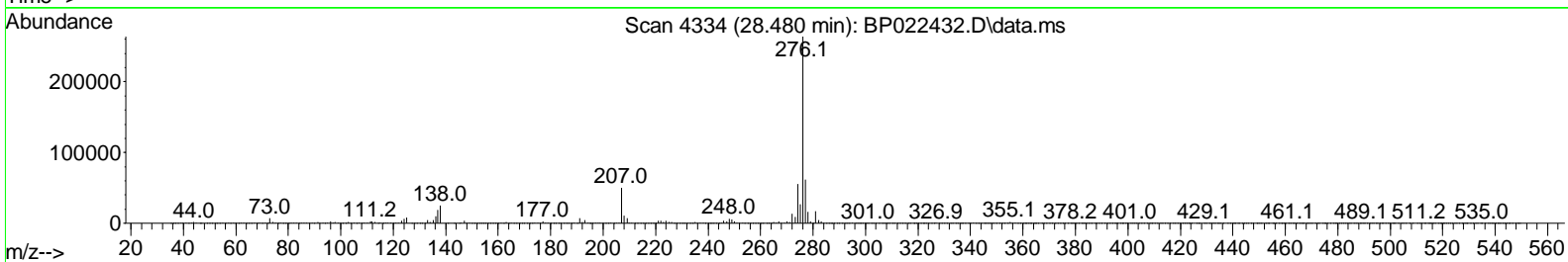
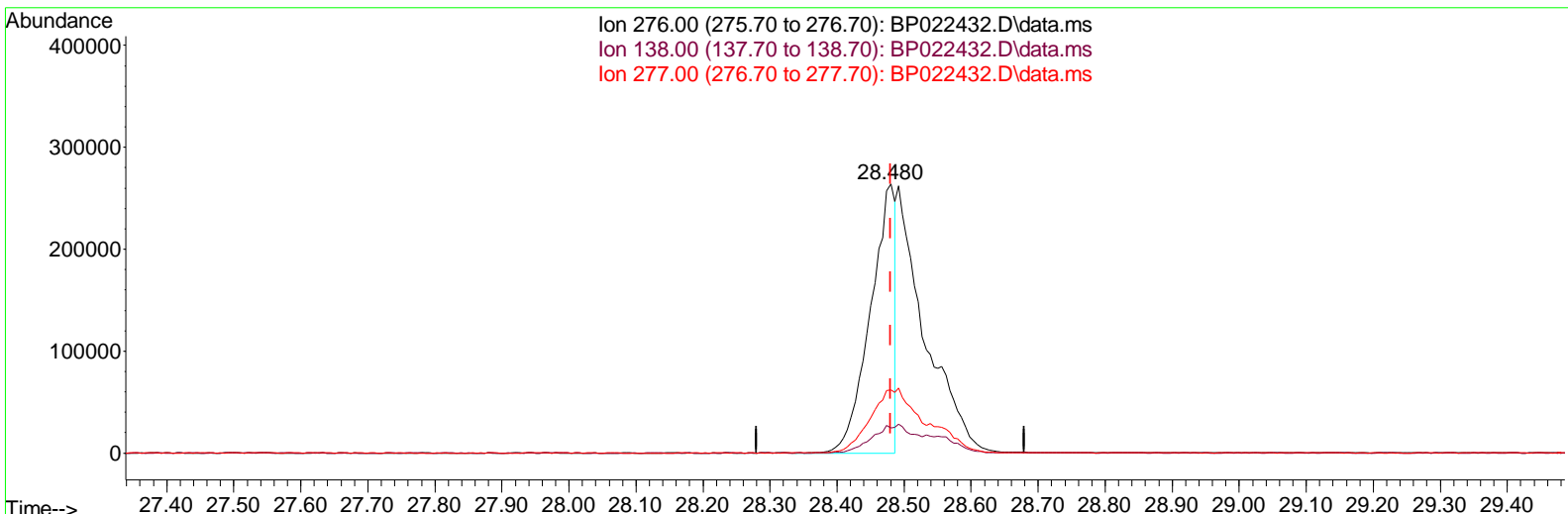
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101624\
 Data File : BP022432.D
 Acq On : 17 Oct 2024 03:17
 Operator : RC/JU
 Sample : SSTDCCC020
 Mi sc :
 ALS Vial : 23 Sample Multi plier: 1

Instrument :
 BNA_P
LabSampleId :
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Oct 17 05:46:30 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 17 05:42:26 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/18/2024
 Supervised By :mohammad ahmed 10/18/2024



TIC: BP022432.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

28.480min (0.000) 8.88 ng/ul

response 676740

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	11.40	9.41
277.00	23.20	23.55
0.00	0.00	0.00

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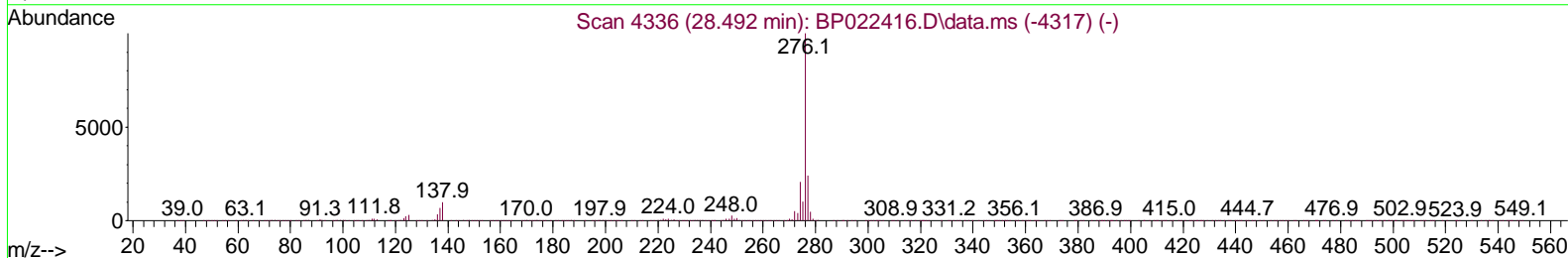
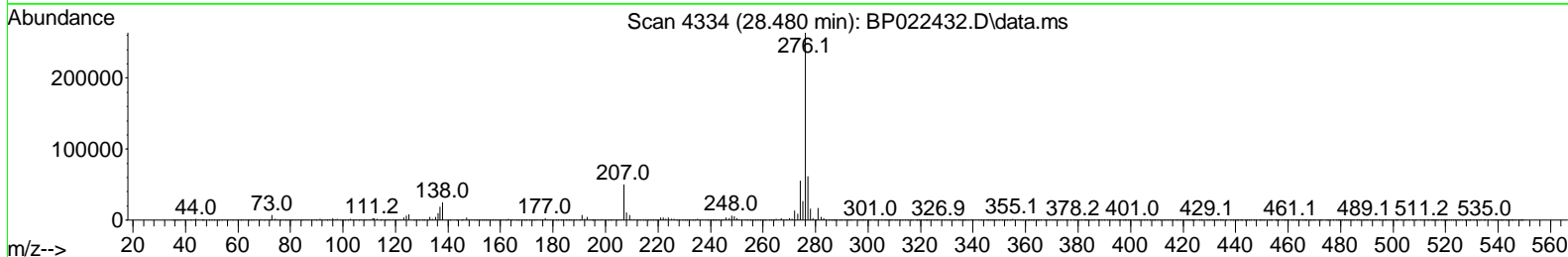
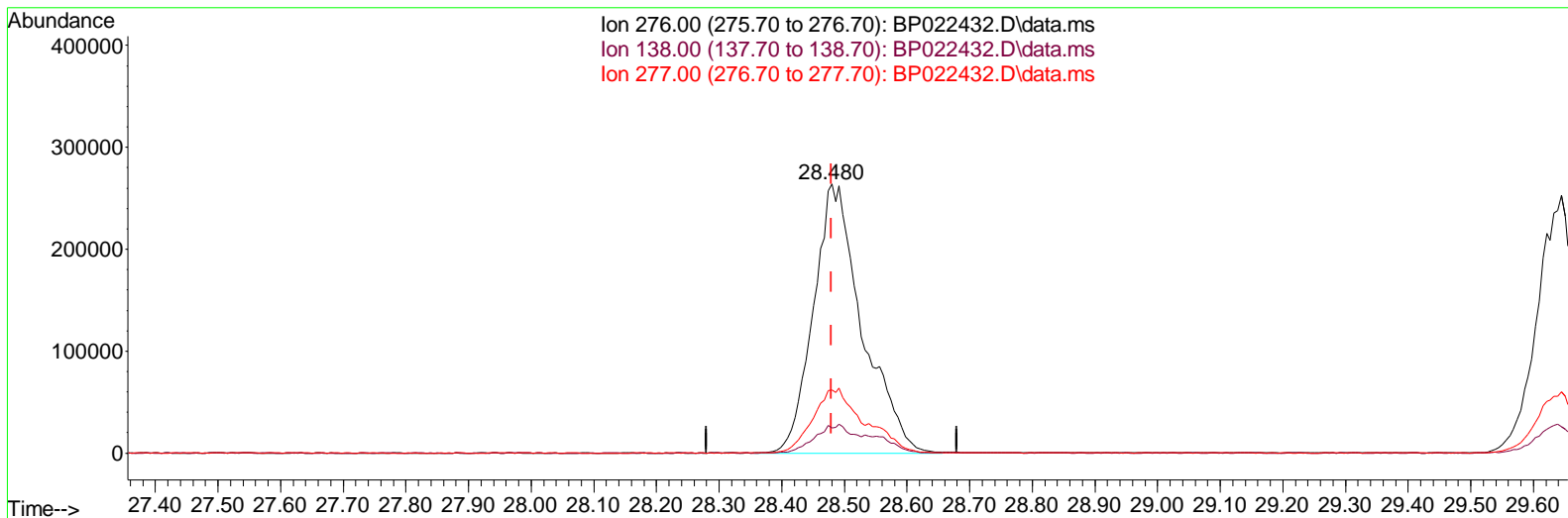
SSTDCCC020

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TIC: BP022432.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

28.480min (0.000) 18.71 ng/ul m

response 1426165

Ion	Exp%	Act%
276.00	100.00	100.00
138.00	11.40	9.41
277.00	23.20	23.55
0.00	0.00	0.00

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Lab Sampled :
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Manual Integrations APPROVED

Reviewed By :Yogesh Patel 10/18/2024
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Quant Time: Oct 17 05:48:59 2024
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 Quant Title : SVOA CALI BRATI ON
 QLast Update : Thu Oct 17 05:42:26 2024
 Response via : Initial Calibration

Compound	R. T.	QI on	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Di chlorobenzene-d4	7.663	152	98817	20.000	ng/ul	0.00
20) Naphthalene-d8	10.422	136	386689	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	319685	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.075	188	823936	20.000	ng/ul	0.00
79) Chrysene-d12	21.516	240	867162	20.000	ng/ul	0.00
88) Perylene-d12	24.780	264	985943	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	17401	6.843	ng/uL	0.00
4) Pyridine-d5	3.611	84	126892	18.177	ng/ul	0.00
7) Phenol-d5	6.858	99	161922	19.466	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.011	67	92754	18.972	ng/ul	0.00
11) 2-Chlorophenol-d4	7.210	132	121708	20.620	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113	132504	19.961	ng/ul	0.00
21) Nitrobenzene-d5	8.805	128	60935	20.494	ng/ul	0.00
24) 2-Nitrophenol-d4	9.522	143	73237	21.276	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.057	165	151440	20.700	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	166735	19.287	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	506547	19.943	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	539843	20.011	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	84759	19.891	ng/ul	0.00
60) Fluorene-d10	15.275	176	441772	20.052	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.422	200	99545	18.370	ng/ul	0.00
73) Anthracene-d10	17.186	188	748874	19.321	ng/ul	0.00
81) Pyrene-d10	19.563	212	949269	20.701	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.551	264	1019378	19.360	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.234	88	19655	7.189	ng/uL	98
5) Pyridine	3.628	79	131786	18.411	ng/ul	100
6) Benzaldehyde	6.828	77	105329	23.432	ng/ul	98
8) Phenol	6.887	94	162667	18.796	ng/ul	97
10) Bis(2-Chloroethyl)ether	7.099	93	128850	19.894	ng/ul	95
12) 2-Chlorophenol	7.240	128	121414	19.893	ng/ul	95
13) 2-Methylphenol	8.105	108	123311	19.634	ng/ul	97
14) 2,2'-oxybis(1-Chloropr...	8.181	45	143967	19.126	ng/ul	98
16) Acetophenone	8.475	105	216239	19.731	ng/ul	95
17) N-Nitrosodimethylamine	8.457	70	112712	20.300	ng/ul	97
18) 4-Methylphenol	8.428	108	136692	19.619	ng/ul	92
19) Hexachloroethane	8.728	117	58003	20.406	ng/ul	98
22) Nitrobenzene	8.846	77	169820	19.135	ng/ul	96
23) Isophorone	9.369	82	322977	20.297	ng/ul	98
25) 2-Nitrophenol	9.552	139	76421	20.600	ng/ul	94
26) 2,4-Dimethylphenol	9.610	107	159880	19.828	ng/ul	99
27) Bis(2-Chloroethoxy)met...	9.840	93	182590	20.209	ng/ul	95
29) 2,4-Dichlorophenol	10.081	162	146729	20.366	ng/ul	99
30) Naphthalene	10.475	128	407094	19.766	ng/ul	98
32) 4-Chloroaniline	10.581	127	165460	19.390	ng/ul	98
33) Hexachlorobutadiene	10.751	225	122120	20.111	ng/ul	100
34) Caprolactam	11.357	113	44886	19.325	ng/ul	92
35) 4-Chloro-3-methylphenol	11.716	107	162348	20.653	ng/ul	99

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Quant Time: Oct 17 05: 48: 59 2024
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 Quant Title : SVOA CALI BRATI ON
 QLast Update : Thu Oct 17 05: 42: 26 2024
 Response via : Ini tial Cal i brati on

Compound	R. T.	QI on	Response	Conc	Units	Dev(Mi n)
36) 2-Methyl naphthal ene	12.081	142	305701	20.220	ng/ul	98
37) 1-Methyl naphthal ene	12.298	142	318368	20.631	ng/ul	100
39) 1, 2, 4, 5-Tetrachl oroben. . .	12.451	216	227232	19.223	ng/ul	97
40) Hexachl orocycl opentadi ene	12.428	237	135029	18.749	ng/ul	98
41) 2, 4, 6-Tri chl orophenol	12.698	196	148547	19.978	ng/ul	98
42) 2, 4, 5-Tri chl orophenol	12.775	196	162153	20.481	ng/ul	97
43) 1, 1' -Bi phenyl	13.098	154	443222	19.235	ng/ul	99
44) 2-Chl oronaphthal ene	13.140	162	368108	19.515	ng/ul	99
45) 2-Ni troani li ne	13.351	65	105314	18.698	ng/ul	91
47) Di methyl phthal ate	13.734	163	504931	19.950	ng/ul	100
48) 2, 6-Di ni trotol uene	13.851	165	101197	21.139	ng/ul	96
50) Acenaphthyl ene	13.993	152	581084	20.754	ng/ul	99
51) 3-Ni troani li ne	14.187	138	92043	20.471	ng/ul	98
52) Acenaphthene	14.340	153	386258	19.509	ng/ul	98
53) 2, 4-Di ni trophenol	14.410	184	73670	20.916	ng/ul	98
55) 4-Ni trophenol	14.516	109	92980	19.660	ng/ul	94
56) Di benzofuran	14.681	168	585445	19.686	ng/ul	97
57) 2, 4-Di ni trotol uene	14.651	165	155489	20.902	ng/ul	97
58) 2, 3, 4, 6-Tetrachl orophenol	14.910	232	157364	21.295	ng/ul	100
59) Di ethyl phthal ate	15.122	149	505947	20.836	ng/ul	98
61) Fl uorene	15.339	166	486470	20.076	ng/ul	98
62) 4-Chl orophenyl -phenyl e. . .	15.334	204	278654	20.109	ng/ul	99
63) 4-Ni troani li ne	15.363	138	101450	22.423	ng/ul	97
66) 4, 6-Di ni tro-2-methyl ph. . .	15.439	198	105706	18.110	ng/ul	99
67) N-Ni trosodi phenyl ami ne	15.557	169	417610	18.927	ng/ul	100
68) 4-Bromophenyl -phenyl ether	16.245	248	203134	20.444	ng/ul	97
69) Hexachl orobenzene	16.363	284	247931	20.272	ng/ul	95
70) Atrazi ne	16.534	200	178287	19.439	ng/ul	93
71) Pentachl orophenol	16.722	266	164403	20.906	ng/ul	99
72) Phenanthrene	17.122	178	848683	19.253	ng/ul	100
74) Anthracene	17.222	178	865083	19.665	ng/ul	98
75) 1, 2, 3, 4-Tetrachl oroben. . .	13.063	216	225228	17.953	ng/uL	98
76) Pentachl orobenzene	14.598	250	260402	18.776	ng/uL	99
77) Carbazol e	17.498	167	746648	19.182	ng/ul	99
78) Di -n-butyl phthal ate	18.086	149	885795	20.567	ng/ul	100
80) Fl uoranthene	19.210	202	1140343	21.631	ng/ul	99
82) Pyrene	19.598	202	1161235	20.552	ng/ul	98
83) Butyl benzyl phthal ate	20.539	149	432091	21.964	ng/ul	96
84) 3, 3' -Di chl orobenzi di ne	21.422	252	425335	20.576	ng/ul	99
85) Benzo(a)anthracene	21.498	228	1207097	19.856	ng/ul	99
86) Bi s(2-ethyl hexyl)phtha. . .	21.427	149	624777	22.126	ng/ul	98
87) Chrysene	21.563	228	1095389	19.433	ng/ul	100
89) Di -n-octyl phthal ate	22.674	149	1046182	22.308	ng/ul	100
90) Benzo(b)fl uoranthene	23.745	252	1251914	20.172	ng/ul #	99
91) Benzo(k)fl uoranthene	23.810	252	1189413	19.575	ng/ul	99
93) Benzo(a)pyrene	24.621	252	1124766	19.692	ng/ul	99
94) I ndeno(1, 2, 3-cd)pyrene	28.480	276	1426165m	18.711	ng/ul	
95) Di benzo(a, h)anthracene	28.556	278	1178515	19.094	ng/ul	98
96) Benzo(g, h, i)peryl ene	29.645	276	1129686	19.054	ng/ul	100

(#) = qual i fier out of range (m) = manual i ntegrati on (+) = signal s summed

Instrument :

BNA_P

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