

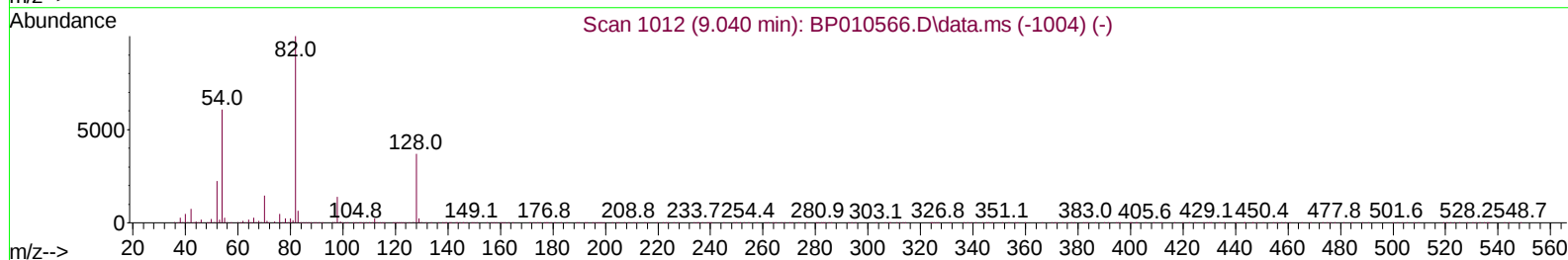
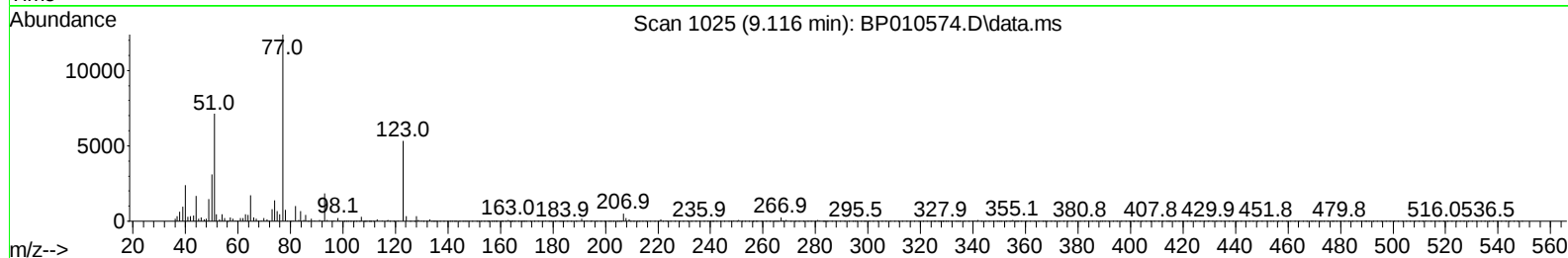
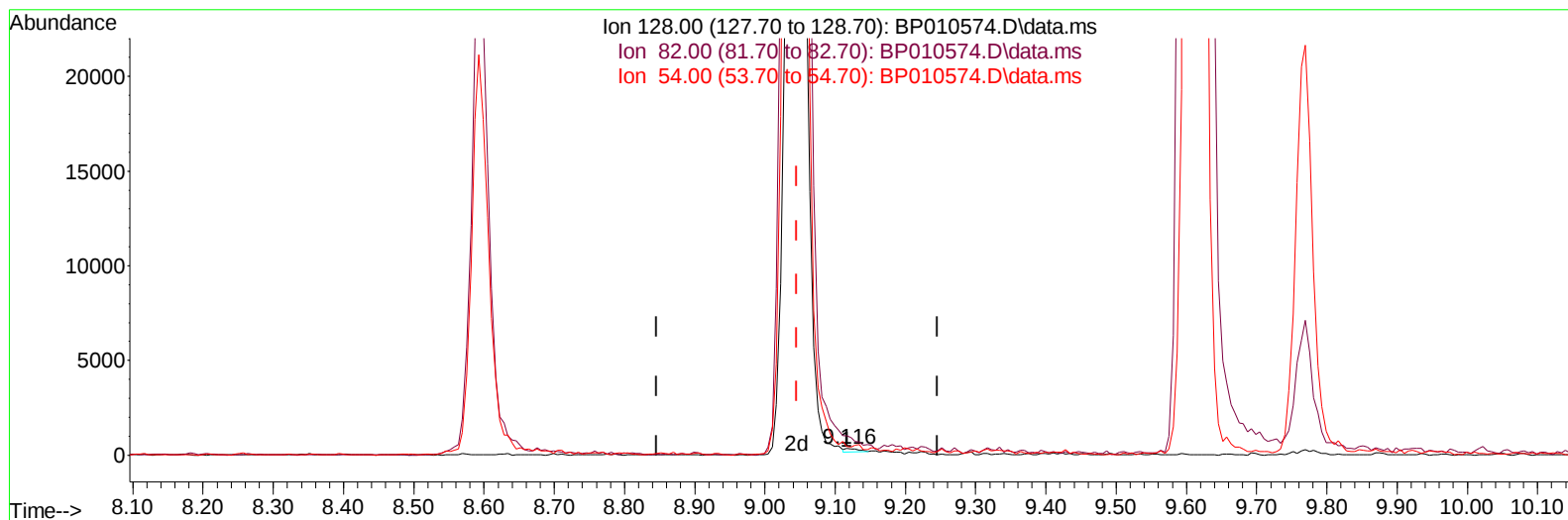
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052722\
 Data File : BP010574.D
 Acq On : 27 May 2022 11:01
 Operator : CG/JU
 Sample : N2815-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 X3A87

Manual IntegrationsAPPROVED

Quant Time: May 27 12:47:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 24 05:50:40 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/27/2022
 Supervised By :mohammad ahmed 05/31/2022



TIC: BP010574.D\data.ms

(21) Nitrobenzene-d5 (S)

9.116min (+ 0.071) 0.06 ng/u1

response 232

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	287.30	291.43
54.00	120.20	127.43
0.00	0.00	0.00

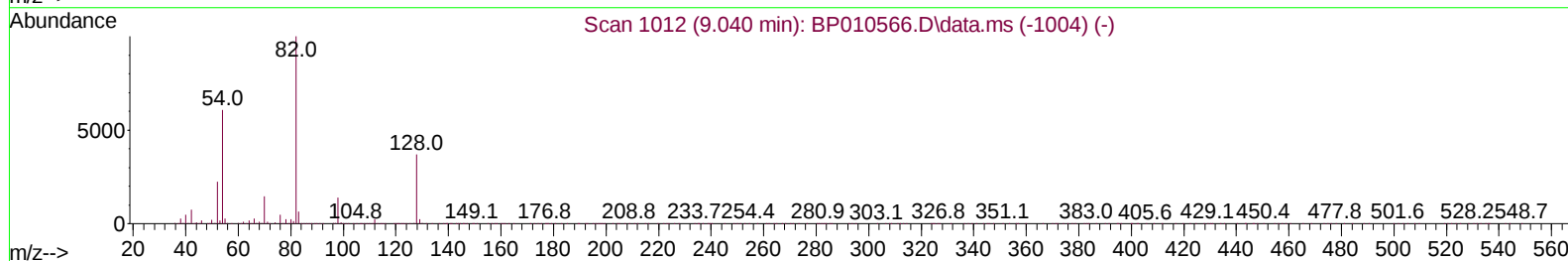
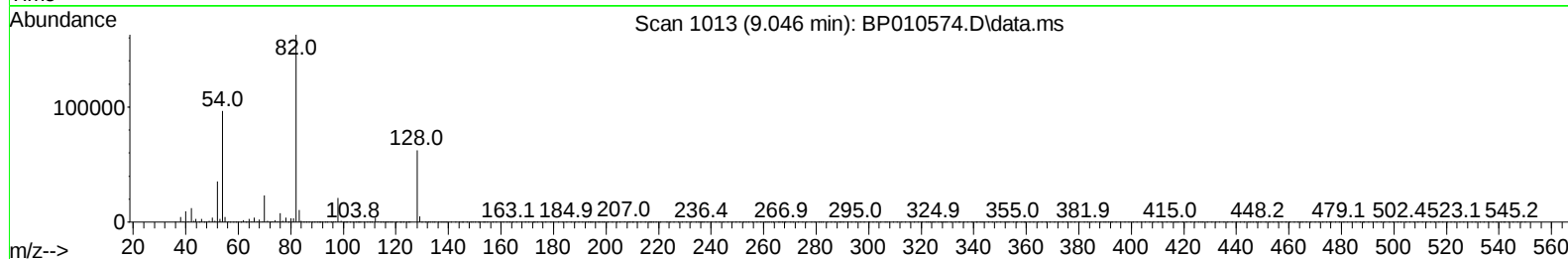
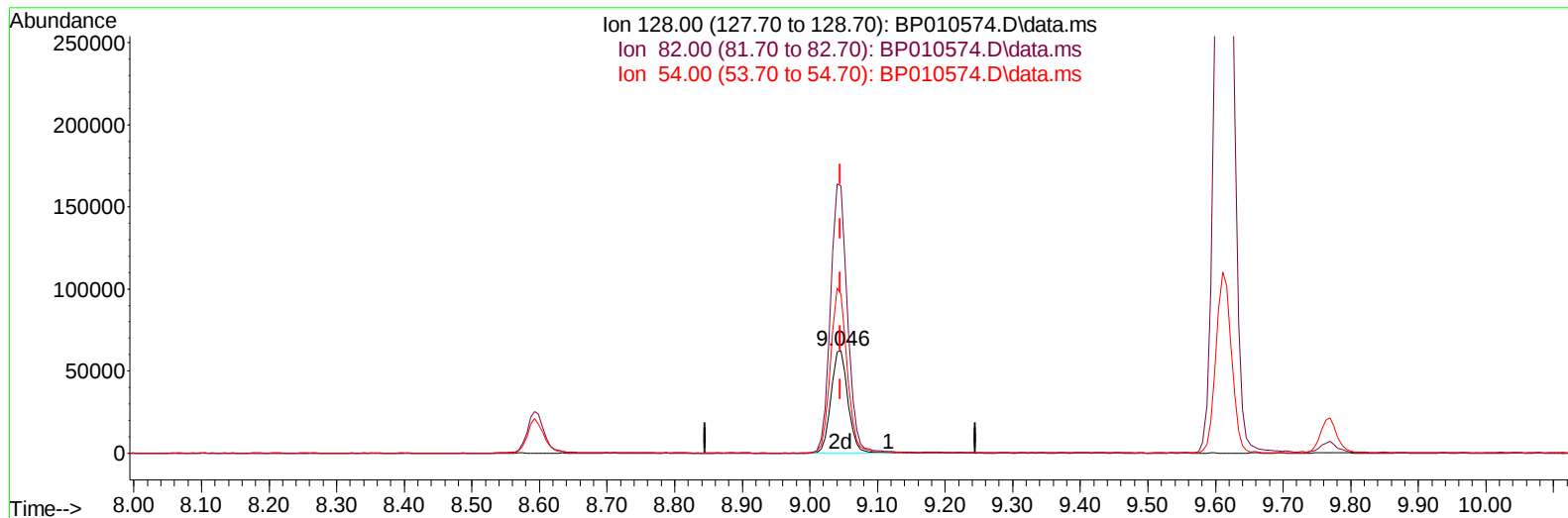
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TIC: BP010574.D\data.ms

(21) Nitrobenzene-d5 (S)

9.046min (+ 0.000) 26.02 ng/ul m

response 108755

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	287.30	260.22
54.00	120.20	154.52#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	122434	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	540045	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.516	164	366781	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.275	188	814328	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	768881	20.000	ng/ul	# 0.00
88) Perylene-d12	23.780	264	668331	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	11888	4.139	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.052	99	253813	24.699	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	175943	25.483	ng/ul	0.00
11) 2-Chlorophenol-d4	7.416	132	200685	25.522	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	201769	23.104	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	108755m	26.020	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	110414	24.928	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	190702	22.770	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	155133	12.457	ng/ul	0.00
46) Dimethylphthalate-d6	13.928	166	763613	28.258	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	818070	26.245	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	97217	20.212	ng/ul	0.00
60) Fluorene-d10	15.516	176	678233	28.849	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	102596	21.542	ng/ul	0.00
73) Anthracene-d10	17.375	188	1231662	33.402	ng/ul	0.00
81) Pyrene-d10	19.604	212	1508927	36.954	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.627	264	1181063	35.193	ng/ul	0.01
Target Compounds						
					Qvalue	
8) Phenol	7.075	94	494465	44.602	ng/ul#	90
22) Nitrobenzene	9.087	77	449105	40.889	ng/ul#	83
23) Isophorone	9.610	82	975701	48.061	ng/ul	97
30) Naphthalene	10.734	128	1569862	55.654	ng/ul	97
36) 2-Methylnaphthalene	12.346	142	874261	43.781	ng/ul	91
41) 2,4,6-Trichlorophenol	12.957	196	335694	48.114	ng/ul#	84
43) 1,1'-Biphenyl	13.351	154	946805	35.169	ng/ul#	93
50) Acenaphthylene	14.240	152	735496	21.812	ng/ul	95
56) Dibenzofuran	14.916	168	1533236	47.561	ng/ul	97
59) Diethylphthalate	15.340	149	1143543	41.809	ng/ul#	93
69) Hexachlorobenzene	16.581	284	353644	36.660	ng/ul#	87
71) Pentachlorophenol	16.928	266	371972	61.454	ng/ul#	84
74) Anthracene	17.410	178	2117606	48.571	ng/ul	98
77) Carbazole	17.681	167	2322228	60.608	ng/ul#	96
80) Fluoranthene	19.281	202	2480313	52.270	ng/ul	97
85) Benzo(a)anthracene	21.345	228	1983288	41.249	ng/ul	95
86) Bis(2-ethylhexyl)phtha...	21.269	149	1621404	55.801	ng/ul#	81
89) Di-n-octyl phthalate	22.198	149	3013289	60.101	ng/ul	100
94) Indeno(1,2,3-cd)pyrene	26.315	276	2045990	52.548	ng/ul#	93
95) Dibenzo(a,h)anthracene	26.327	278	1790578	52.999	ng/ul#	95
96) Benzo(g,h,i)perylene	27.086	276	1802350	55.437	ng/ul#	90

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev	(Min)

(#) = qualifier out of range (m) = manual integration (+) = signals summed								

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