

Quantitation Report (Qedit)

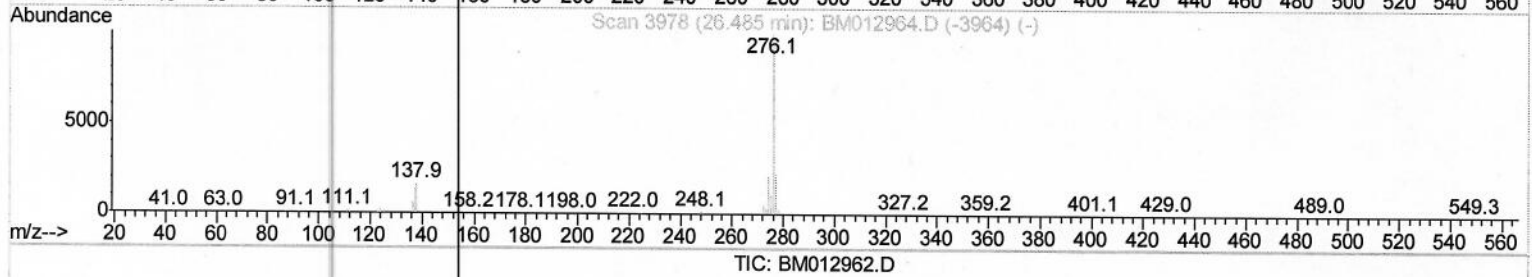
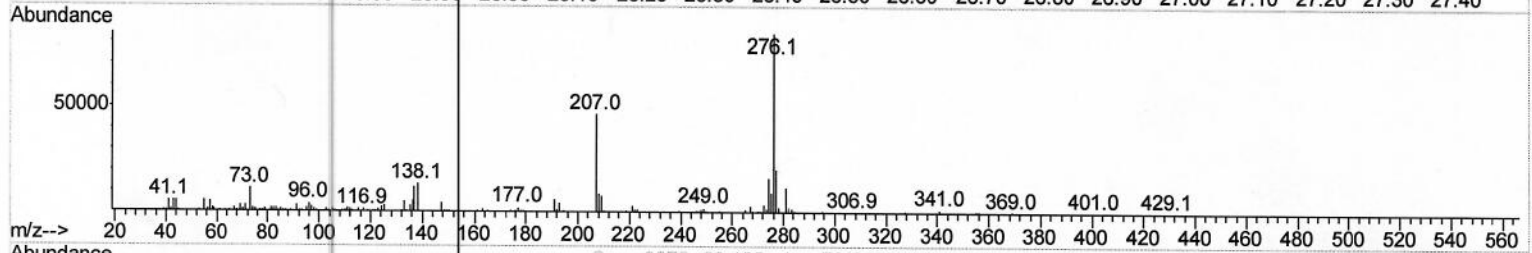
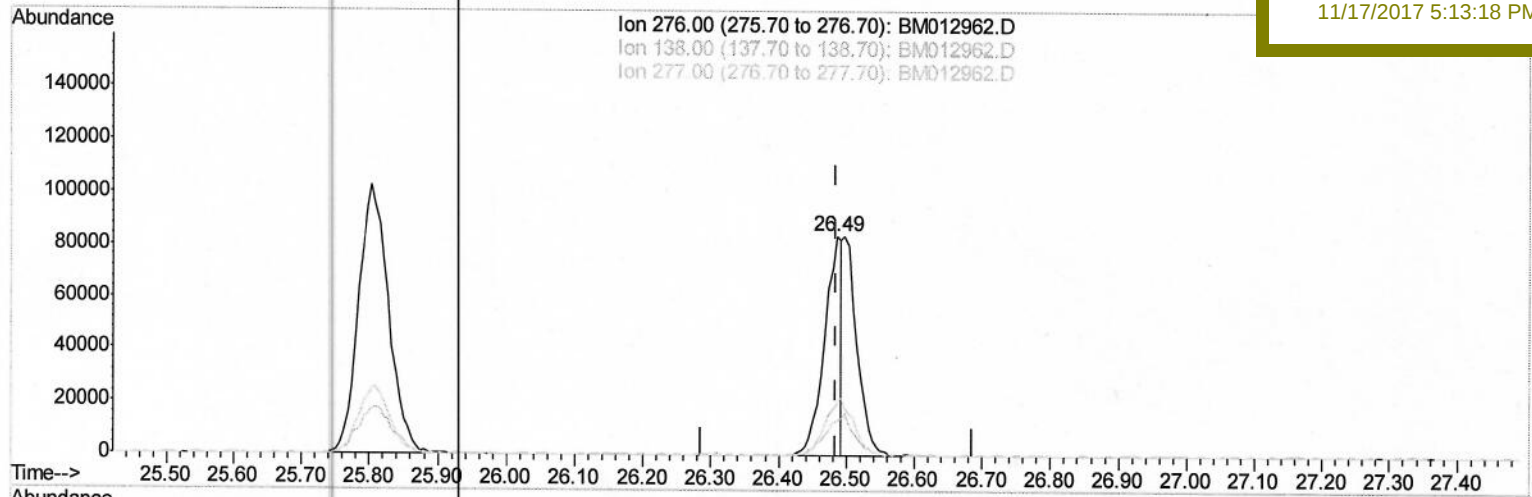
Data Path : Z:\HPCHEM1\BNA M\Data\BM111617\
 Data File : BM012962.D
 Acq On : 16 Nov 2017 13:10
 Operator : SJ/JU
 Sample : SSTD00526
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Nov 16 16:13:38 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111617.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 16 16:08:46 2017
 Response via : Initial Calibration

Instrument :
 BNA_M
ClientSampleId :
 SSTD00526

Manual Integrations
APPROVED

SOHIL
 11/17/2017 5:13:18 PM



(91) Benzo(g,h,i)perylene
 26.485min (+0.000) 2.97ng/ul
 response 148215

Ion	Exp%	Act%
276.00	100	100
138.00	10.60	15.89#
277.00	23.30	24.13
0.00	0.00	0.00

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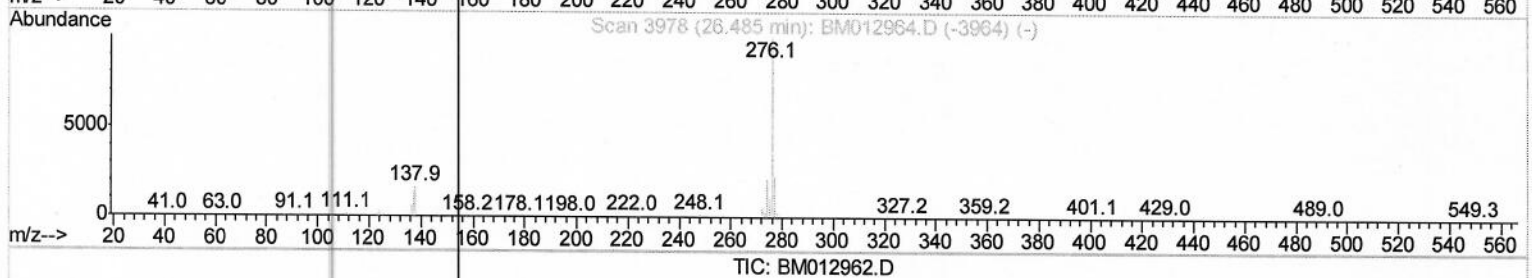
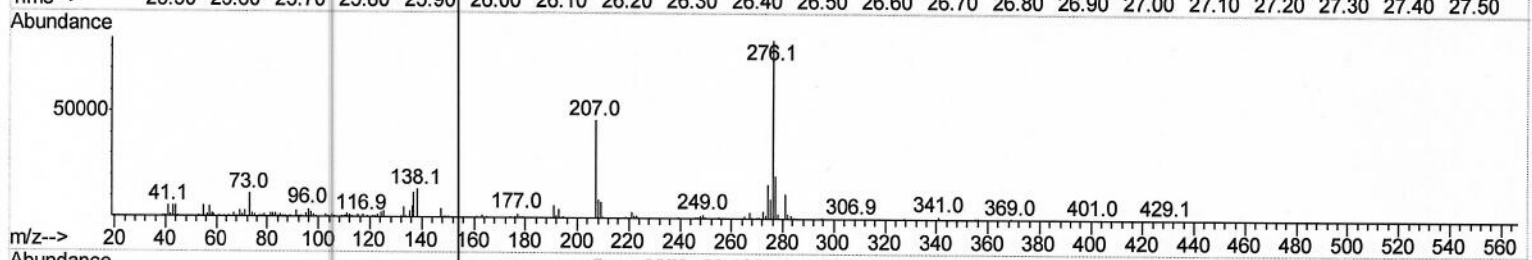
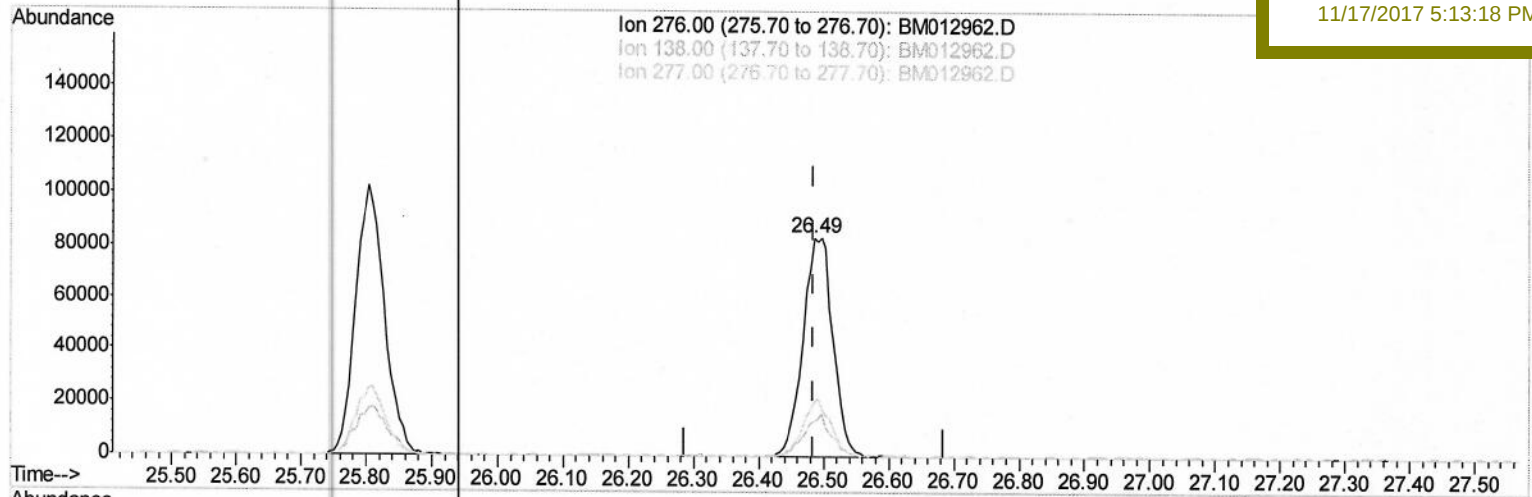
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11/17/2017 5:13:18 PM



(91) Benzo(g,h,i)perylene

26.485min (+0.000) 5.32ng/ul m > SJ 11/21/17

response 265294

Ion	Exp%	Act%
276.00	100	100
138.00	10.60	15.89#
277.00	23.30	24.13
0.00	0.00	0.00

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	155298	20.00	ng/ul	0.00
18) Naphthalene-d8	10.52	136	620919	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	371353	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.12	188	834890	20.00	ng/ul	0.00
75) Chrysene-d12	21.30	240	905374	20.00	ng/ul	0.00
83) Perylene-d12	23.54	264	920710	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.26	96	7856	2.19	ng/uL	0.00
5) Phenol-d5	0.00	99	0d	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0d	0.00	ng/ul	
9) 2-Chlorophenol-d4	7.26	132	50979	4.46	ng/ul	0.00
13) 4-Methylphenol-d8	0.00	113	0d	0.00	ng/ul	
19) Nitrobenzene-d5	8.88	128	20726	4.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.60	143	19133	3.68	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	54243	5.33	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
43) Dimethylphthalate-d6	13.79	166	164649	5.03	ng/ul	0.00
46) Acenaphthylene-d8	14.06	160	206786	5.11	ng/ul	0.00
51) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
57) Fluorene-d10	15.37	176	138210	5.07	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
70) Anthracene-d10	17.22	188	214004	5.06	ng/ul	0.00
76) Pyrene-d10	19.51	212	232609	5.17	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.40	264	226790	4.93	ng/ul	0.00

Target Compounds

				Ovalue	
2) 1,4-Dioxane	3.29	88	7705	1.882	ng/uL# 59
10) 2-Chlorophenol	7.29	128	51221	4.493	ng/ul 99
15) N-Nitroso-di-n-propylamine	8.53	70	46029	4.840	ng/ul# 67
17) Hexachloroethane	8.81	117	21416	4.776	ng/ul# 73
20) Nitrobenzene	8.92	77	54190	4.547	ng/ul 95
21) Isophorone	9.45	82	119180	5.313	ng/ul# 93
23) 2-Nitrophenol	9.63	139	20607	3.669	ng/ul# 75
24) 2,4-Dimethylphenol	9.69	107	63271	5.397	ng/ul# 88
25) Bis(2-Chloroethoxy)methane	9.94	93	65668	4.699	ng/ul 91
27) 2,4-Dichlorophenol	10.16	162	48672	5.087	ng/ul# 92
28) Naphthalene	10.56	128	162527	4.917	ng/ul 99
31) Hexachlorobutadiene	10.86	225	35822	5.966	ng/ul# 89
33) 4-Chloro-3-methylphenol	11.80	107	53231	4.929	ng/ul 98
34) 2-Methylnaphthalene	12.19	142	119057	4.960	ng/ul 91
36) 1,2,4,5-Tetrachlorobenzene	12.55	216	66696	5.492	ng/ul# 97
38) 2,4,6-Trichlorophenol	12.79	196	38424	4.990	ng/ul 94
39) 2,4,5-Trichlorophenol	12.86	196	41084	4.992	ng/ul 95
40) 1,1'-Biphenyl	13.20	154	154314	4.860	ng/ul# 98
41) 2-Chloronaphthalene	13.25	162	123901	5.081	ng/ul 96
42) 2-Nitroaniline	13.44	65	21631	3.060	ng/ul 94
44) Dimethylphthalate	13.84	163	155275	4.917	ng/ul 100
45) 2,6-Dinitrotoluene	13.95	165	19782	3.296	ng/ul# 88
47) Acenaphthylene	14.09	152	190439	4.940	ng/ul 100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
49) Acenaphthene	14.43	153	125886	4.678	ng/ul	97
53) Dibenzofuran	14.77	168	183040	4.906	ng/ul	97
54) 2,4-Dinitrotoluene	14.73	165	24779	2.724	ng/ul#	81
55) 2,3,4,6-Tetrachlorophenol	15.00	232	33883	4.593	ng/ul#	91
56) Diethylphthalate	15.21	149	156190	4.874	ng/ul	97
58) Fluorene	15.43	166	149631	4.876	ng/ul	95
59) 4-Chlorophenyl-phenylether	15.43	204	78333	5.085	ng/ul	94
64) N-Nitrosodiphenylamine	15.64	169	131037	4.968	ng/ul	98
65) 4-Bromophenyl-phenylether	16.32	248	49367	5.039	ng/ul	90
66) Hexachlorobenzene	16.42	284	54674	4.820	ng/ul#	90
69) Phenanthrene	17.16	178	229482	4.735	ng/ul	99
71) Anthracene	17.25	178	245399	4.953	ng/ul	97
73) Di-n-butylphthalate	18.11	149	269389	5.178	ng/ul#	97
77) Pyrene	19.54	202	280427	4.991	ng/ul#	90
78) Butylbenzylphthalate	20.46	149	114787	5.038	ng/ul	96
80) Benzo(a)anthracene	21.29	228	294552	5.250	ng/ul	99
81) Bis(2-ethylhexyl)phthalate	21.26	149	169743	5.130	ng/ul#	99
82) Chrysene	21.34	228	272300	5.175	ng/ul	99
85) Benzo(b)fluoranthene	22.87	252	289782	5.010	ng/ul#	96
86) Benzo(k)fluoranthene	22.92	252	268194	4.798	ng/ul#	98
88) Benzo(a)pyrene	23.45	252	274965	5.051	ng/ul#	98
89) Indeno(1,2,3-cd)pyrene	25.80	276	312947	5.123	ng/ul#	95
90) Dibenzo(a,h)anthracene	25.83	278	267838	5.209	ng/ul#	96
91) Benzo(a,h,i)perylene	26.49	276	265294m)	5.321	ng/ul	96

5511/21/17

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

(QT Reviewed)

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