

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061118\
 Data File : BM015635.D
 Acq On : 12 Jun 2018 11:40
 Operator : SJ/JU
 Sample : J3346-01
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DAXG3

Quant Time: Jun 12 14:12:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM060818MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 12 04:48:05 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.34	152	155908	20.00	ng/ul	0.00
18) Naphthalene-d8	11.17	136	603787	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.96	164	295431	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.68	188	592454	20.00	ng/ul	0.00
77) Chrysene-d12	21.79	240	611132	20.00	ng/ul	0.00
85) Perylene-d12	24.20	264	668522	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.55	96	15878	4.95	ng/uL	0.00
5) Phenol-d5	7.49	99	284403	20.09	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.65	67	191838	22.43	ng/ul	0.00
9) 2-Chlorophenol-d4	7.86	132	252672	23.35	ng/ul	0.00
13) 4-Methylphenol-d8	9.05	113	207989	17.06	ng/ul	0.00
19) Nitrobenzene-d5	9.52	128	111934	27.57	ng/ul	0.00
22) 2-Nitrophenol-d4	10.25	143	126405	28.95	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.79	165	215524	23.84	ng/ul	0.00
29) 4-Chloroaniline-d4	11.31	131	164877	16.36	ng/ul	0.00
43) Dimethylphthalate-d6	14.36	166	600311	24.54	ng/ul	0.00
46) Acenaphthylene-d8	14.66	160	766742	29.07	ng/ul	0.00
51) 4-Nitrophenol-d4	15.16	143	85264	18.06	ng/ul	0.01
57) Fluorene-d10	15.94	176	503887	24.35	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.07	200	56859	15.93	ng/ul	0.00
70) Anthracene-d10	17.77	188	723150	26.77	ng/ul	0.00
78) Pyrene-d10	20.03	212	753174	28.65	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.04	264	885607	26.43	ng/ul	0.00

Target Compounds

					Qvalue	
6) Phenol	7.52	94	32266	2.195	ng/ul#	86
44) Dimethylphthalate	14.40	163	234761	9.400	ng/ul	99
76) Fluoranthene	19.70	202	78563	1.971	ng/ul	98
79) Pyrene	20.06	202	83288	2.429	ng/ul	99
82) Benzo(a)anthracene	21.77	228	48649	1.313	ng/ul#	89
84) Chrysene	21.82	228	62952	1.811	ng/ul#	84
87) Benzo(b)fluoranthene	23.46	252	132384	3.332	ng/ul#	84
88) Benzo(k)fluoranthene	23.46	252	132384	3.531	ng/ul#	83
90) Benzo(a)pyrene	24.09	252	62881	1.656	ng/ul#	78
91) Indeno(1,2,3-cd)pyrene	26.69	276	58785	1.236	ng/ul	93
93) Benzo(g,h,i)perylene	27.46	276	47455	1.211	ng/ul#	92

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

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