

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121315\
 Data File : BM003508.D
 Acq On : 13 Dec 2015 13:58
 Operator : UM/IZ
 Sample : G4734-16
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4KX7

Quant Time: Dec 14 04:26:25 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM120115.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 14 04:23:11 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	99800	20.00	ng/ul	0.00
18) Naphthalene-d8	10.51	136	454941	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.37	164	242566	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.12	188	492295	20.00	ng/ul	0.00
78) Chrysene-d12	21.31	240	514159	20.00	ng/ul	0.00
86) Perylene-d12	23.57	264	508815	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	4664	2.32	ng/uL	0.00
5) Phenol-d5	6.89	99	89779	11.27	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	140979	32.07	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	191918	29.13	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	136906	20.48	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	100053	32.24	ng/ul	0.00
22) 2-Nitrophenol-d4	9.60	143	108734	33.52	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	185372	28.82	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	18166	2.50	ng/ul	0.00
44) Dimethylphthalate-d6	13.78	166	548672	30.09	ng/ul	0.00
47) Acenaphthylene-d8	14.06	160	670195	29.90	ng/ul	0.00
52) 4-Nitrophenol-d4	14.57	143	23799	7.28	ng/ul	0.00
58) Fluorene-d10	15.36	176	473113	30.15	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.49	200	64649	25.67	ng/ul	0.00
71) Anthracene-d10	17.22	188	670966	30.68	ng/ul	0.00
79) Pyrene-d10	19.52	212	721156	28.65	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.42	264	693671	28.26	ng/ul	0.00

Target Compounds

						Ovalue
6) Phenol	6.92	94	34003	4.04	ng/ul	99
28) Naphthalene	10.57	128	2839944	123.63	ng/ul	99

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

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