

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM121115\
 Data File : BM003483.D
 Acq On : 11 Dec 2015 15:17
 Operator : UM/IZ
 Sample : PB87218BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK18

Quant Time: Dec 12 00:44:06 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM120115.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 12 00:41:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	73579	20.00	ng/ul	0.00
18) Naphthalene-d8	10.51	136	358577	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.37	164	219490	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.12	188	467029	20.00	ng/ul	0.00
78) Chrysene-d12	21.32	240	379165	20.00	ng/ul	0.00
86) Perylene-d12	23.57	264	344443	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	1940	1.31	ng/uL	0.00
5) Phenol-d5	6.89	99	37917	6.45	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	87594	27.03	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	111468	22.95	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	70626	14.33	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	68946	28.19	ng/ul	0.00
22) 2-Nitrophenol-d4	9.60	143	73479	28.74	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	130034	25.65	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	0.00
44) Dimethylphthalate-d6	13.78	166	496646	30.10	ng/ul	0.00
47) Acenaphthylene-d8	14.06	160	579700	28.58	ng/ul	0.00
52) 4-Nitrophenol-d4	14.57	143	10626	3.59	ng/ul	0.00
58) Fluorene-d10	15.37	176	441392	31.08	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.49	200	44627	18.68	ng/ul	0.00
71) Anthracene-d10	17.22	188	684114	32.97	ng/ul	0.00
79) Pyrene-d10	19.52	212	676601	36.45	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.43	264	545273	32.81	ng/ul	0.00

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

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