

Data Path : Z:\HPCHEM1\BNA_M\Data\BM080215\
 Data File : BM002166.D
 Acq On : 02 Aug 2015 12:28
 Operator : TP
 Sample : PB84595BL
 Misc : SBLK95
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK95

Quant Time: Aug 03 01:44:20 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 03 01:33:08 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	88256	20.00	ng/ul	0.00
18) Naphthalene-d8	10.35	136	399979	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.23	164	262866	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.98	188	627868	20.00	ng/ul	0.00
78) Chrysene-d12	21.19	240	791373	20.00	ng/ul	0.00
86) Perylene-d12	23.37	264	817912	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.00	96	11501	6.74	ng/uL	0.00
5) Phenol-d5	6.74	99	268024	42.48	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.90	67	151190	43.33	ng/ul	0.00
9) 2-Chlorophenol-d4	7.09	132	226862	42.42	ng/ul	0.00
13) 4-Methylphenol-d8	8.27	113	239491	47.75	ng/ul	0.00
19) Nitrobenzene-d5	8.72	128	122644	43.14	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	145653	46.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.97	165	252299	41.47	ng/ul	0.00
29) 4-Chloroaniline-d4	10.49	131	351280	52.54	ng/ul	0.00
44) Dimethylphthalate-d6	13.64	166	884616	45.94	ng/ul	0.00
47) Acenaphthylene-d8	13.92	160	1033803	42.94	ng/ul	0.00
52) 4-Nitrophenol-d4	14.46	143	148235	46.12	ng/ul	0.00
58) Fluorene-d10	15.23	176	750566	44.30	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.37	200	161910	42.16	ng/ul	0.00
71) Anthracene-d10	17.08	188	1250616	44.62	ng/ul	0.00
79) Pyrene-d10	19.39	212	1459748	43.81	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.24	264	1762275	44.45	ng/ul	0.01

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

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