

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080616\
 Data File : BM006945.D
 Acq On : 06 Aug 2016 19:33
 Operator : UM/SJ
 Sample : PB92639BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK39

Quant Time: Aug 08 03:52:48 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 08 03:51:46 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	145367	20.00	ng/ul	0.00
7) Naphthalene-d8	10.33	136	727123	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.21	164	496996	20.00	ng/ul	0.00
23) Phenanthrene-d10	16.97	188	1244749	20.00	ng/ul	0.00
29) Chrysene-d12	21.17	240	1374850	20.00	ng/ul	0.00
34) Perylene-d12	23.36	264	1107198	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.09	96	11126	3.50	ng/uL	0.00
3) Phenol-d5	6.77	99	410600	29.79	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.90	67	321090	37.15	ng/ul	0.00
5) 2-Chlorophenol-d4	7.10	132	345530	35.12	ng/ul	0.00
6) 4-Methylphenol-d8	8.29	113	378658	32.91	ng/ul	0.00
8) Nitrobenzene-d5	8.72	128	183383	33.96	ng/ul	0.00
9) 2-Nitrophenol-d4	9.43	143	221469	34.24	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	9.98	165	346253	28.24	ng/ul	0.00
12) 4-Chloroaniline-d4	10.49	131	550216	37.59	ng/ul	0.00
16) Dimethylphthalate-d6	13.63	166	1639950	37.50	ng/ul	0.00
17) Acenaphthylene-d8	13.90	160	1832600	36.23	ng/ul	0.00
20) 4-Nitrophenol-d4	14.50	143	184195	29.03	ng/ul	0.01
21) Fluorene-d10	15.21	176	1356657	36.49	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.37	200	100233	19.50	ng/ul	0.00
26) Anthracene-d10	17.06	188	2226313	37.02	ng/ul	0.00
30) Pyrene-d10	19.37	212	2557401	39.02	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.22	264	1999162	37.91	ng/ul	0.00

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

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