

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072116\
 Data File : BM006586.D
 Acq On : 21 Jul 2016 14:16
 Operator : UM/SJ
 Sample : PB92235BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK35

Quant Time: Jul 21 15:42:11 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 21 01:09:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	142434	20.00	ng/ul	0.00
7) Naphthalene-d8	10.39	136	609691	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.26	164	343567	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.02	188	779303	20.00	ng/ul	0.00
29) Chrysene-d12	21.22	240	916000	20.00	ng/ul	0.00
34) Perylene-d12	23.41	264	945013	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 1,4-Dioxane-d8	3.13	96	13835	3.93	ng/uL	0.00
3) Phenol-d5	6.79	99	496588	40.98	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.95	67	317764	42.33	ng/ul	0.00
5) 2-Chlorophenol-d4	7.15	132	394679	40.76	ng/ul	0.00
6) 4-Methylphenol-d8	8.32	113	389922	41.49	ng/ul	0.00
8) Nitrobenzene-d5	8.77	128	191817	41.56	ng/ul	0.00
9) 2-Nitrophenol-d4	9.48	143	210831	42.96	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.02	165	365000	39.36	ng/ul	0.00
12) 4-Chloroaniline-d4	10.53	131	537527	52.63	ng/ul	0.00
16) Dimethylphthalate-d6	13.68	166	1113631	41.34	ng/ul	0.00
17) Acenaphthylene-d8	13.96	160	1418494	41.26	ng/ul	0.00
20) 4-Nitrophenol-d4	14.48	143	177129	37.86	ng/ul	0.00
21) Fluorene-d10	15.26	176	977437	40.57	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.40	200	141493	34.98	ng/ul	0.00
26) Anthracene-d10	17.11	188	1512168	40.87	ng/ul	0.00
30) Pyrene-d10	19.42	212	1688701	40.46	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.27	264	1778934	40.92	ng/ul	0.00

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

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