

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021916\
 Data File : BM004305.D
 Acq On : 20 Feb 2016 04:07
 Operator : SJ/UM
 Sample : PB88454BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK54

Quant Time: Feb 20 06:05:57 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM021216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Feb 20 06:01:29 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	20989	20.00	ng/ul	0.00
18) Naphthalene-d8	10.42	136	110256	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	75587	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.04	188	187127	20.00	ng/ul	0.00
78) Chrysene-d12	21.26	240	230978	20.00	ng/ul	0.00
86) Perylene-d12	23.48	264	231338	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.09	96	2520	6.05	ng/uL	0.00
5) Phenol-d5	6.82	99	55332	32.82	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.96	67	31662	34.77	ng/ul	0.00
9) 2-Chlorophenol-d4	7.17	132	49270	32.95	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	53321	37.13	ng/ul	0.00
19) Nitrobenzene-d5	8.79	128	26928	31.76	ng/ul	0.00
22) 2-Nitrophenol-d4	9.51	143	31631	33.38	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	51518	30.15	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	87160	42.84	ng/ul	0.00
44) Dimethylphthalate-d6	13.70	166	230175	36.59	ng/ul	0.00
47) Acenaphthylene-d8	13.98	160	260059	34.03	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	32253	31.69	ng/ul	0.00
58) Fluorene-d10	15.29	176	193609	36.52	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.44	200	30600	26.85	ng/ul	0.00
71) Anthracene-d10	17.14	188	321160	35.10	ng/ul	0.00
79) Pyrene-d10	19.45	212	387768	31.63	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.34	264	428712	34.73	ng/ul	0.00

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

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