

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050416\
 Data File : BM005227.D
 Acq On : 04 May 2016 18:04
 Operator : UM/SJ
 Sample : PB89256BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK56

Quant Time: May 04 18:41:35 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM050416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 04 17:24:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	8372	20.00	ng/ul	0.00
18) Naphthalene-d8	10.57	136	42868	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.42	164	30522	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.17	188	89447	20.00	ng/ul	0.00
75) Chrysene-d12	21.36	240	148960	20.00	ng/ul	0.00
83) Perylene-d12	23.64	264	178640	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.27	96	923	14.62	ng/uL	-0.01
5) Phenol-d5	6.95	99	20169	23.84	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.12	67	13606	30.47	ng/ul	0.00
9) 2-Chlorophenol-d4	7.32	132	16157	28.39	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	17182	24.47	ng/ul	0.00
19) Nitrobenzene-d5	8.94	128	7601	28.97	ng/ul	0.00
22) 2-Nitrophenol-d4	9.66	143	7982	27.41	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	15033	25.20	ng/ul	0.00
29) 4-Chloroaniline-d4	10.71	131	26258	31.81	ng/ul	0.00
43) Dimethylphthalate-d6	13.83	166	79787	32.25	ng/ul	0.00
46) Acenaphthylene-d8	14.11	160	82758	29.69	ng/ul	0.00
51) 4-Nitrophenol-d4	14.63	143	10385	20.73	ng/ul	0.00
57) Fluorene-d10	15.42	176	71250	31.87	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.54	200	8064	16.30	ng/ul	0.00
70) Anthracene-d10	17.26	188	130828	31.95	ng/ul	0.00
76) Pyrene-d10	19.56	212	181442	30.78	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.49	264	255201	32.39	ng/ul	0.00

Target Compounds

					Ovalue
21) Isophorone	9.66	82	568	0.40 ng/ul#	65
45) 2,6-Dinitrotoluene	13.83	165	391	0.91 ng/ul#	33
88) Benzo(a)pyrene	23.49	252	418	0.04 ng/ul#	60

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

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