

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022616\
 Data File : BM004426.D
 Acq On : 26 Feb 2016 13:08
 Operator : SJ/UM
 Sample : PB88533BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK33

Quant Time: Feb 26 23:25:18 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM022216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 26 23:20:20 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	20679	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	93147	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.28	164	54322	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	112767	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	111397	20.00	ng/ul	0.00
86) Perylene-d12	23.47	264	109323	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	2686	6.85	ng/uL	0.00
5) Phenol-d5	6.80	99	50514	29.13	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	27918	31.66	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	47283	31.62	ng/ul	0.00
13) 4-Methylphenol-d8	8.34	113	44986	29.03	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	22863	32.26	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	25759	31.89	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	42760	29.28	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	65682	36.09	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	152097	32.65	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	184308	33.79	ng/ul	0.00
52) 4-Nitrophenol-d4	14.53	143	14856	19.84	ng/ul	0.01
58) Fluorene-d10	15.28	176	129355	32.65	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.43	200	14337	20.70	ng/ul	0.00
71) Anthracene-d10	17.13	188	190948	34.56	ng/ul	0.00
79) Pyrene-d10	19.44	212	198538	36.84	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.33	264	204436	35.63	ng/ul	0.00

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

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