

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM022616\
 Data File : BM004423.D
 Acq On : 26 Feb 2016 10:25
 Operator : SJ/UM
 Sample : PB88557BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK57

Quant Time: Feb 26 23:24:38 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	7.62	152	20766	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	100465	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.28	164	63112	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	135458	20.00	ng/ul	0.00
78) Chrysene-d12	21.25	240	114142	20.00	ng/ul	0.00
86) Perylene-d12	23.47	264	105728	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	2701	6.86	ng/uL	0.00
5) Phenol-d5	6.80	99	55630	31.95	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	30956	34.96	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	51811	34.50	ng/ul	0.00
13) 4-Methylphenol-d8	8.34	113	51156	32.88	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	26341	34.46	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	30373	34.86	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	50148	31.84	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	78758	40.12	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	191801	35.44	ng/ul	0.00
47) Acenaphthylene-d8	13.97	160	228272	36.02	ng/ul	0.00
52) 4-Nitrophenol-d4	14.53	143	20527	23.59	ng/ul	0.00
58) Fluorene-d10	15.28	176	161769	35.14	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.43	200	20074	24.12	ng/ul	0.00
71) Anthracene-d10	17.13	188	239901	36.14	ng/ul	0.00
79) Pyrene-d10	19.44	212	235110	42.58	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.33	264	207967	37.48	ng/ul	0.00

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

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