

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021516\
 Data File : BM004279.D
 Acq On : 16 Feb 2016 04:58
 Operator : SJ/UM
 Sample : H1416-22
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9QF5

Quant Time: Feb 16 05:54:38 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM021216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 16 02:45:44 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	33914	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	157648	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.31	164	87289	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.06	188	170781	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	150541	20.00	ng/ul	0.01
86) Perylene-d12	23.50	264	150938	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.11	96	973	1.45	ng/uL	0.00
5) Phenol-d5	6.85	99	24420	8.96	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	6.99	67	49125	33.39	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	73726	30.52	ng/ul	0.00
13) 4-Methylphenol-d8	8.38	113	47591	20.51	ng/ul	0.02
19) Nitrobenzene-d5	8.81	128	42756	35.27	ng/ul	0.00
22) 2-Nitrophenol-d4	9.53	143	48695	35.94	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.07	165	78452	32.11	ng/ul	0.01
29) 4-Chloroaniline-d4	10.59	131	3587	1.23	ng/ul	0.02
44) Dimethylphthalate-d6	13.72	166	247984	34.14	ng/ul	0.00
47) Acenaphthylene-d8	14.00	160	305348	34.60	ng/ul	0.00
52) 4-Nitrophenol-d4	14.58	143	7312	6.22	ng/ul	0.04
58) Fluorene-d10	15.30	176	205282	33.53	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	28754	27.65	ng/ul	0.00
71) Anthracene-d10	17.16	188	289605	34.69	ng/ul	0.00
79) Pyrene-d10	19.47	212	277652	34.75	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	272544	33.84	ng/ul	0.01

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

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