

Data Path : Z:\HPCHEM1\BNA_M\Data\BM121416\
 Data File : BM008302.D
 Acq On : 14 Dec 2016 09:36
 Operator : UM/SJ
 Sample : H5963-10
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SVOC-GPC2-BLANK

Quant Time: Dec 14 11:06:06 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 14 05:32:31 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	109656	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	419691	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.32	164	282104	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.07	188	565852	20.00	ng/ul	0.00
75) Chrysene-d12	21.27	240	745375	20.00	ng/ul	0.00
83) Perylene-d12	23.50	264	760926	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0	0.00	ng/ul	
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	10.46	131	246	0.03	ng/ul	-0.15
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
46) Acenaphthylene-d8	14.03	160	277	0.01	ng/ul	0.02
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
57) Fluorene-d10	15.36	176	1575	0.10	ng/ul	0.05
62) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
70) Anthracene-d10	17.07	188	565852	22.95	ng/ul	-0.09
76) Pyrene-d10	19.50	212	1136	0.04	ng/ul	0.03
87) Benzo(a)pyrene-d12	23.50	264	719256	22.51	ng/ul	0.15

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

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