

Data Path : Z:\HPCHEM1\BNA G\DATA\BG112116\
 Data File : BG024876.D
 Acq On : 22 Nov 2016 2:18
 Operator : UM/SJ
 Sample : H5694-18
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JHSS3

Manual Integrations
 APPROVED

sohil
 11/22/2016 4:07:43 PM

Quant Time: Nov 22 06:32:49 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG111416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 22 01:29:33 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.26	152	126416	20.00	ng/ul	0.00
18) Naphthalene-d8	11.09	136	522858	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.88	164	443560	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.62	188	956666m	20.00	ng/ul	0.00
75) Chrysene-d12	21.92	240	1157826	20.00	ng/ul	0.00
83) Perylene-d12	25.36	264	1238208	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.63	96	10147	4.16	ng/uL	0.00
5) Phenol-d5	7.39	99	207449	19.15	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.57	67	146780	21.54	ng/ul	0.00
9) 2-Chlorophenol-d4	7.78	132	159399	20.52	ng/ul	0.00
13) 4-Methylphenol-d8	8.95	113	186708	20.53	ng/ul	0.00
19) Nitrobenzene-d5	9.43	128	91367	23.08	ng/ul	0.00
22) 2-Nitrophenol-d4	10.16	143	99724	21.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.70	165	194657	20.69	ng/ul	0.00
29) 4-Chloroaniline-d4	11.22	131	205890	19.59	ng/ul	0.00
43) Dimethylphthalate-d6	14.27	166	654973	20.25	ng/ul	0.00
46) Acenaphthylene-d8	14.57	160	798595	22.66	ng/ul	0.00
51) 4-Nitrophenol-d4	15.06	143	80950	14.07	ng/ul	0.00
57) Fluorene-d10	15.86	176	636924	20.92	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.97	200	67133m	10.21	ng/ul	0.00
70) Anthracene-d10	17.72	188	968386	22.92	ng/ul	0.00
76) Pyrene-d10	19.99	212	1130729	22.52	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.11	264	1216523	22.44	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	7.42	94	28132	2.49	ng/ul 95
44) Dimethylphthalate	14.31	163	110567	3.56	ng/ul 98

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

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