

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101316\
 Data File : BG024342.D
 Acq On : 14 Oct 2016 5:03
 Operator : UM/SJ
 Sample : H5187-22
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BDPE3

Manual Integrations
 APPROVED

Umangi
 10/14/2016 6:43:36 PM

Quant Time: Oct 14 11:47:33 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG0100716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 14 06:45:57 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	214864	20.00	ng/ul	0.00
18) Naphthalene-d8	11.12	136	932157	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.90	164	716487	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.64	188	1399271	20.00	ng/ul	0.00
75) Chrysene-d12	21.94	240	1581446	20.00	ng/ul	0.00
83) Perylene-d12	25.38	264	1643139	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.65	96	5210	1.15	ng/uL	0.00
5) Phenol-d5	7.41	99	104596	5.22	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.60	67	282726	20.86	ng/ul	0.00
9) 2-Chlorophenol-d4	7.81	132	251704	18.23	ng/ul	0.00
13) 4-Methylphenol-d8	8.97	113	193577	11.94	ng/ul	0.00
19) Nitrobenzene-d5	9.46	128	160458	24.26	ng/ul	0.00
22) 2-Nitrophenol-d4	10.19	143	190585	25.75	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.72	165	333584	20.69	ng/ul	0.00
29) 4-Chloroaniline-d4	11.24	131	6255m	0.37	ng/ul	0.00
43) Dimethylphthalate-d6	14.29	166	1208082	24.15	ng/ul	0.00
46) Acenaphthylene-d8	14.60	160	1372817	23.96	ng/ul	0.00
51) 4-Nitrophenol-d4	15.06	143	49899	5.63	ng/ul	0.00
57) Fluorene-d10	15.89	176	1126395	23.81	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.99	200	240267	28.57	ng/ul	0.00
70) Anthracene-d10	17.74	188	1723495	27.27	ng/ul	0.00
76) Pyrene-d10	20.01	212	1934179	26.20	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.15	264	2024642	27.39	ng/ul	0.00

Target Compounds Ovalue

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

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