

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102517\
 Data File : BG029456.D
 Acq On : 25 Oct 2017 23:57
 Operator : SJ/JU
 Sample : I5793-18
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B0DA3

Quant Time: Oct 26 07:18:32 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG101417.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 24 23:55:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	76320	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.99	136	364013	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.78	164	241371	20.00	ng/ul	-0.01
61) Phenanthrene-d10	17.53	188	600691	20.00	ng/ul	0.00
75) Chrysene-d12	21.79	240	663937	20.00	ng/ul	-0.01
83) Perylene-d12	25.10	264	669139	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.55	96	12591	6.71	ng/uL	0.00
5) Phenol-d5	7.31	99	246355	33.63	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	7.48	67	149773	32.67	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.70	132	196941	38.11	ng/ul	-0.01
13) 4-Methylphenol-d8	8.86	113	219867	37.16	ng/ul	-0.01
19) Nitrobenzene-d5	9.33	128	103298	39.53	ng/ul	-0.01
22) 2-Nitrophenol-d4	10.06	143	118529	44.27	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.60	165	223129	36.57	ng/ul	-0.01
29) 4-Chloroaniline-d4	11.11	131	34728	4.90	ng/ul	-0.01
43) Dimethylphthalate-d6	14.18	166	730803	35.41	ng/ul	0.00
46) Acenaphthylene-d8	14.48	160	910308	36.27	ng/ul	0.00
51) 4-Nitrophenol-d4	14.96	143	117524	30.85	ng/ul	-0.01
57) Fluorene-d10	15.77	176	675880	39.99	ng/ul	-0.01
62) 4,6-Dinitro-2-methylphenol	15.88	200	114125	38.92	ng/ul	-0.01
70) Anthracene-d10	17.62	188	1057804	37.23	ng/ul	-0.01
76) Pyrene-d10	19.89	212	1243522	36.18	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.88	264	1224839	38.65	ng/ul	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
6) Phenol	7.34	94	28747	3.793	ng/ul	88

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

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