

Data File : BG027485.D
Acq On : 17 Jun 2017 00:01
Operator : SJ/MA
Sample : I3493-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BDC59

Quant Time: Jun 17 03:25:55 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 16 22:56:07 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.52	152	37868	20.00	ng/ul	0.00
18) Naphthalene-d8	11.34	136	188978	20.00	ng/ul	0.00
35) Acenaphthene-d10	15.10	164	127218	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.85	188	324359	20.00	ng/ul	0.00
75) Chrysene-d12	22.23	240	330480	20.00	ng/ul	0.00
83) Perylene-d12	25.89	264	315357	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	4.08	96	5848	6.18	ng/uL	0.00
5) Phenol-d5	7.65	99	26389	6.43	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.83	67	89049	31.81	ng/ul	0.00
9) 2-Chlorophenol-d4	8.05	132	68030	25.07	ng/ul	0.00
13) 4-Methylphenol-d8	9.19	113	47449	14.78	ng/ul	0.00
19) Nitrobenzene-d5	9.68	128	43643	31.79	ng/ul	0.00
22) 2-Nitrophenol-d4	10.40	143	47667	32.68	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.94	165	78907	28.64	ng/ul	0.00
29) 4-Chloroaniline-d4	11.46	131	5090	1.46	ng/ul	0.00
43) Dimethylphthalate-d6	14.49	166	339958	31.92	ng/ul	0.00
46) Acenaphthylene-d8	14.80	160	396063	32.28	ng/ul	0.00
51) 4-Nitrophenol-d4	15.27	143	11927	5.72	ng/ul	0.00
57) Fluorene-d10	16.09	176	302434	30.27	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.19	200	55944	27.77	ng/ul	0.00
70) Anthracene-d10	17.95	188	488995	32.33	ng/ul	0.00
76) Pyrene-d10	20.22	212	556523	34.48	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.63	264	472961	33.31	ng/ul	0.00

Target Compounds Qvalue

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

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