

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062718\
 Data File : BG035467.D
 Acq On : 28 Jun 2018 5:56
 Operator : JU/SJ
 Sample : J3711-02DL 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 QNWP8-27S-GWDL

Quant Time: Jun 28 07:46:46 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 08 17:16:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	48014	20.00	ng	-0.02
21) Naphthalene-d8	10.85	136	187367	20.00	ng	-0.03
38) Acenaphthene-d10	14.67	164	136777	20.00	ng	-0.02
63) Phenanthrene-d10	17.42	188	360419	20.00	ng	-0.02
75) Chrysene-d12	21.68	240	405676	20.00	ng	-0.04
86) Perylene-d12	24.90	264	394882	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	78547	27.61	ng	-0.02
7) Phenol-d6	7.21	99	69693	16.60	ng	-0.01
23) Nitrobenzene-d5	9.21	82	169517	43.01	ng	-0.03
41) 2,4,6-Tribromophenol	16.16	330	117425	67.06	ng	-0.02
44) 2-Fluorobiphenyl	13.29	172	386169	36.70	ng	-0.04
78) Terphenyl-d14	20.02	244	547549	28.27	ng	-0.03
Target Compounds						
2) 1,4-Dioxane	3.55	88	3520	2.593	ng	Qvalue # 58
27) 2,4-Dimethylphenol	10.02	122	31247	10.685	ng	84
31) Naphthalene	10.91	128	457351	46.987	ng	100
49) Dimethylphthalate	14.12	163	30741	2.577	ng	99
51) Acenaphthene	14.74	154	67642	7.425	ng	97
54) Dibenzofuran	15.07	168	68159	4.996	ng	91
57) Fluorene	15.72	166	50948	4.468	ng	93
70) Phenanthrene	17.46	178	77597	4.238	ng	98
72) Carbazole	17.82	167	127587	7.498	ng	99

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

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