

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062618\
 Data File : BG035444.D
 Acq On : 27 Jun 2018 13:40
 Operator : JU/SJ
 Sample : J3711-02
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 QNWP8-27S-GW

Quant Time: Jun 27 16:07:17 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	48588	20.00	ng	-0.02
21) Naphthalene-d8	10.85	136	186453	20.00	ng	-0.03
38) Acenaphthene-d10	14.67	164	133413	20.00	ng	-0.02
63) Phenanthrene-d10	17.42	188	339426	20.00	ng	-0.02
75) Chrysene-d12	21.68	240	376692	20.00	ng	-0.04
86) Perylene-d12	24.90	264	364942	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	154425	53.64	ng	-0.01
7) Phenol-d6	7.21	99	147898	34.80	ng	-0.01
23) Nitrobenzene-d5	9.22	82	347211	88.52	ng	-0.02
41) 2,4,6-Tribromophenol	16.16	330	248489	145.50	ng	-0.02
44) 2-Fluorobiphenyl	13.30	172	764819	74.53	ng	-0.03
78) Terphenyl-d14	20.02	244	1028944	57.22	ng	-0.02
Target Compounds						
2) 1,4-Dioxane	3.55	88	5872	4.275	ng	Qvalue # 72
10) Phenol	7.24	94	9151	2.081	ng	94
27) 2,4-Dimethylphenol	10.02	122	69197	23.778	ng	98
31) Naphthalene	10.91	128	913749	94.337	ng	99
45) 1,1'-Biphenyl	13.50	154	40420	3.768	ng	93
49) Dimethylphthalate	14.12	163	63857	5.489	ng	98
51) Acenaphthene	14.74	154	144919	16.309	ng	97
54) Dibenzofuran	15.07	168	144727	10.875	ng	97
57) Fluorene	15.72	166	99221	8.921	ng	97
70) Phenanthrene	17.46	178	156642	9.085	ng	99
72) Carbazole	17.82	167	242171	15.112	ng	98

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

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