

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090518\
 Data File : BG036567.D
 Acq On : 5 Sep 2018 19:17
 Operator : JU/SJ
 Sample : J4741-28
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A0C2-04-D

Quant Time: Sep 06 02:57:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 31 13:03:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	57703	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	245428	20.00	ng	0.00
38) Acenaphthene-d10	14.68	164	168534	20.00	ng	0.00
63) Phenanthrene-d10	17.42	188	424140	20.00	ng	0.00
75) Chrysene-d12	21.69	240	443828	20.00	ng	0.00
86) Perylene-d12	24.90	264	443723	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	177546	48.81	ng	0.00
7) Phenol-d6	7.21	99	207194	39.78	ng	-0.01
23) Nitrobenzene-d5	9.22	82	370027	81.80	ng	0.00
41) 2,4,6-Tribromophenol	16.17	330	267063	132.80	ng	0.00
44) 2-Fluorobiphenyl	13.31	172	883482	74.85	ng	0.00
78) Terphenyl-d14	20.03	244	1501184	79.06	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.91	128	441042	35.949	ng	99
37) 2-Methylnaphthalene	12.51	142	52725	5.873	ng	95
49) Dimethylphthalate	14.13	163	157281	11.429	ng	99
51) Acenaphthene	14.75	154	25345	2.371	ng	90
57) Fluorene	15.73	166	26399	2.109	ng	93
70) Phenanthrene	17.47	178	55964	2.597	ng	95

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

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