

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122217\
 Data File : BG031754.D
 Acq On : 23 Dec 2017 18:19
 Operator : SJ/JU
 Sample : I7037-02
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 VNJ-221-OILY-STONES

Manual Integrations
 APPROVED

Sohil
 12/26/2017 3:31:14 PM

Quant Time: Dec 26 00:40:13 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 11:08:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.83	152	47544	20.00	ng	-0.01
21) Naphthalene-d8	11.71	136	200607	20.00	ng	-0.01
38) Acenaphthene-d10	15.46	164	136148	20.00	ng	-0.01
63) Phenanthrene-d10	18.19	188	337414	20.00	ng	-0.01
75) Chrysene-d12	22.55	240	399516	20.00	ng	-0.02
86) Perylene-d12	26.34	264	428108	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	6.24	112	401756	153.91	ng	0.00
7) Phenol-d6	7.92	99	458828	135.39	ng	0.00
23) Nitrobenzene-d5	10.03	82	307191	115.63	ng	-0.01
41) 2,4,6-Tribromophenol	16.93	330	334048	160.80	ng	-0.01
44) 2-Fluorobiphenyl	14.09	172	1076475	99.24	ng	-0.01
78) Terphenyl-d14	20.73	244	1918698	109.72	ng	-0.01
Target Compounds						
31) Naphthalene	11.77	128	1003521	99.121	ng	99
37) 2-Methylnaphthalene	13.32	142	58579m	7.137	ng	
54) Dibenzofuran	15.85	168	36679	2.575	ng	94
57) Fluorene	16.50	166	33273	2.875	ng	90
70) Phenanthrene	18.23	178	54929	2.877	ng	99
72) Carbazole	18.58	167	41704	2.446	ng	99

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

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