

Data Path : Z:\HPCHEM1\BNA G\DATA\BG102017\
 Data File : BG029389.D
 Acq On : 21 Oct 2017 13:24
 Operator : SJ/JU
 Sample : I5981-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NB-01-101817-A

Manual Integrations
 APPROVED

Sohil
 10/23/2017 3:41:15 PM

Quant Time: Oct 21 23:05:43 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 17:32:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	54479	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	234259	20.00	ng	0.00
38) Acenaphthene-d10	14.78	164	161166	20.00	ng	0.00
63) Phenanthrene-d10	17.52	188	442758	20.00	ng	0.00
75) Chrysene-d12	21.79	240	503587	20.00	ng	-0.02
86) Perylene-d12	25.09	264	511901	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	416793	127.81	ng	0.00
7) Phenol-d6	7.31	99	496169	108.39	ng	0.00
23) Nitrobenzene-d5	9.33	82	351619	95.38	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	353130	169.71	ng	0.00
44) 2-Fluorobiphenyl	13.41	172	961123	87.46	ng	0.00
78) Terphenyl-d14	20.11	244	1910335	86.30	ng	0.00
Target Compounds						
31) Naphthalene	11.03	128	939559	80.476	ng	98
37) 2-Methylnaphthalene	12.63	142	67690	7.955	ng	95
51) Acenaphthene	14.84	154	32076	3.117	ng	95
54) Dibenzofuran	15.18	168	64601	4.398	ng	98
57) Fluorene	15.82	166	36225	2.985	ng	97
70) Phenanthrene	17.56	178	60556	2.647	ng	96

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

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