

Data Path : \\74.0.250.170\TERASTORAGE\SVOASRV\HPCHEM1\BNA_G\DATA\BG062919\
 Data File : BG041547.D
 Acq On : 29 Jun 2019 15:31
 Operator : HP/JU
 Sample : K3568-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RW-1

Manual Integrations
 APPROVED

mohammad
 7/1/2019 5:58:24 PM

Quant Time: Jun 30 01:24:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG062619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 27 13:25:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	29641	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	100173	20.00	ng	0.00
39) Acenaphthene-d10	14.68	164	72864	20.00	ng	0.00
64) Phenanthrene-d10	17.43	188	191607	20.00	ng	0.00
76) Chrysene-d12	21.70	240	232488	20.00	ng	0.00
87) Perylene-d12	24.99	264	257522	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	87588	52.68	ng	0.00
7) Phenol-d6	7.18	99	73491	31.40	ng	0.00
23) Nitrobenzene-d5	9.21	82	214667	89.38	ng	0.00
42) 2,4,6-Tribromophenol	16.17	330	168729	135.62	ng	0.00
45) 2-Fluorobiphenyl	13.30	172	492278	86.70	ng	0.00
79) Terphenyl-d14	20.03	244	1018560	93.10	ng	0.00
Target Compounds						
37) 2-Methylnaphthalene	12.51	142	122593	29.954	ng	98
38) 1-Methylnaphthalene	12.73	142	281774m	72.165	ng	
52) Acenaphthene	14.74	154	17431	3.982	ng	92
55) Dibenzofuran	15.08	168	24417	3.204	ng	# 95
58) Fluorene	15.72	166	35133	5.796	ng	# 99
71) Phenanthrene	17.47	178	40388	3.826	ng	99

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

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