

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120618\
 Data File : BG038390.D
 Acq On : 6 Dec 2018 19:51
 Operator : JU/SJ
 Sample : J6003-23DL 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NON-UST-03R-CDL

Manual Integrations
 APPROVED

mohammad
 12/7/2018 12:20:56 PM

Quant Time: Dec 06 21:38:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG120418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 04 17:15:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	46993	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	125959	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	82171	20.00	ng	0.00
64) Phenanthrene-d10	17.45	188	237492	20.00	ng	0.00
76) Chrysene-d12	21.73	240	214332	20.00	ng	-0.01
87) Perylene-d12	25.04	264	221535	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	64531	24.30	ng	0.00
7) Phenol-d6	7.21	99	69748	18.17	ng	-0.01
23) Nitrobenzene-d5	9.25	82	90865	35.82	ng	0.00
42) 2,4,6-Tribromophenol	16.20	330	58504	58.00	ng	0.00
45) 2-Fluorobiphenyl	13.33	172	218478	37.89	ng	0.00
79) Terphenyl-d14	20.05	244	400660	40.05	ng	0.00
Target Compounds						
31) Naphthalene	10.94	128	29717	4.939	ng	98
37) 2-Methylnaphthalene	12.54	142	156678	36.566	ng	98
38) 1-Methylnaphthalene	12.76	142	162403m	39.573	ng	
58) Fluorene	15.75	166	17328	2.956	ng	98
71) Phenanthrene	17.49	178	39234	3.299	ng	99

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

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