

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122118\
 Data File : BG038812.D
 Acq On : 22 Dec 2018 11:18
 Operator : JU/SJ
 Sample : J6505-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-59(0-23)-COMP

Manual Integrations
 APPROVED

Sohil
 12/26/2018 3:17:38 PM

Quant Time: Dec 24 04:31:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	27418	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	115948	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	79240	20.00	ng	0.00
64) Phenanthrene-d10	17.45	188	174150	20.00	ng	0.00
76) Chrysene-d12	21.73	240	162615	20.00	ng	0.00
87) Perylene-d12	25.03	264	180109	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	178848	115.70	ng	0.01
7) Phenol-d6	7.22	99	255573	120.43	ng	0.02
23) Nitrobenzene-d5	9.24	82	159419	70.24	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	137179	125.61	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	428676	74.21	ng	0.00
79) Terphenyl-d14	20.05	244	613461	82.10	ng	0.00
Target Compounds						
10) Phenol	7.25	94	16151	7.164	ng	87
50) Dimethylphthalate	14.14	163	43393	6.706	ng	# 95
75) Fluoranthene	19.50	202	59603	5.334	ng	98
78) Pyrene	19.86	202	56624	5.271	ng	96
81) Benzo(a)anthracene	21.71	228	31169	3.146	ng	98
83) Chrysene	21.78	228	29565m	3.098	ng	
84) Bis(2-ethylhexyl)phthalate	21.57	149	68990	11.590	ng	99
88) Benzo(b)fluoranthene	23.98	252	40303	4.076	ng	96
90) Benzo(a)pyrene	24.87	252	29767	3.120	ng	96

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

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