

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061819\
 Data File : BG041300.D
 Acq On : 18 Jun 2019 16:51
 Operator : HP/JU
 Sample : K3351-07DL 10X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RT-3860DL

Manual Integrations
 APPROVED

mohammad
 6/21/2019 9:56:50 PM

Quant Time: Jun 18 17:31:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 14:58:02 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	34669	20.00	ng	0.00
21) Naphthalene-d8	10.90	136	130403	20.00	ng	0.00
39) Acenaphthene-d10	14.71	164	97076	20.00	ng	0.00
64) Phenanthrene-d10	17.46	188	242982	20.00	ng	0.00
76) Chrysene-d12	21.74	240	258475	20.00	ng	0.00
87) Perylene-d12	25.05	264	271945	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	25335	13.43	ng	0.00
7) Phenol-d6	7.23	99	36081	13.19	ng	0.00
23) Nitrobenzene-d5	9.26	82	25919	8.36	ng	0.00
42) 2,4,6-Tribromophenol	16.20	330	16677	11.19	ng	0.00
45) 2-Fluorobiphenyl	13.34	172	49244	6.89	ng	0.00
79) Terphenyl-d14	20.06	244	95291	7.31	ng	0.00
Target Compounds						
52) Acenaphthene	14.78	154	66011	11.725	ng	92
55) Dibenzofuran	15.11	168	57744	5.899	ng	98
58) Fluorene	15.76	166	78074	10.139	ng	100
71) Phenanthrene	17.50	178	479900	37.032	ng	99
72) Anthracene	17.60	178	41026	3.211	ng	98
75) Fluoranthene	19.51	202	664652	38.606	ng	98
78) Pyrene	19.87	202	462953	26.558	ng	99
81) Benzo(a)anthracene	21.72	228	133709	7.640	ng	97
83) Chrysene	21.78	228	114903	6.827	ng	98
88) Benzo(b)fluoranthene	23.98	252	70012m	4.154	ng	
90) Benzo(a)pyrene	24.89	252	37431	2.352	ng	# 95

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

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