

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110719\
 Data File : BF117708.D
 Acq On : 7 Nov 2019 16:53
 Operator : JU
 Sample : K5723-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 7-A

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:31:34 PM

Quant Time: Nov 08 02:23:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 17:27:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	316568	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	1224921	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	645038	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	1152870	20.00	ng	0.00
76) Chrysene-d12	14.12	240	730055	20.00	ng	-0.01
87) Perylene-d12	15.63	264	839928	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1674811	95.77	ng	0.00
7) Phenol-d6	6.55	99	2115555	98.11	ng	0.00
23) Nitrobenzene-d5	7.49	82	1266957	61.07	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	578080	91.61	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	2378595	72.76	ng	0.00
79) Terphenyl-d14	13.06	244	2448851	69.05	ng	0.00
Target Compounds						
10) Phenol	6.56	94	108614	4.674	ng	87
50) Dimethylphthalate	9.69	163	425855	9.766	ng	99
71) Phenanthrene	11.49	178	509690	9.224	ng	96
75) Fluoranthene	12.69	202	248271	4.400	ng	95
78) Pyrene	12.92	202	250886	4.708	ng	97
88) Benzo(b)fluoranthene	15.17	252	106443m	2.012	ng	

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

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