

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110518\
 Data File : BF110473.D
 Acq On : 5 Nov 2018 18:57
 Operator : JU/SJ
 Sample : J5812-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP1-A

Manual Integrations
 APPROVED

Sohil
 11/6/2018 9:01:15 AM

Quant Time: Nov 06 04:20:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 16:06:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	117602	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	509713	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	247991	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	443953	20.00	ng	0.00
76) Chrysene-d12	14.13	240	237358	20.00	ng	0.00
87) Perylene-d12	15.66	264	205910	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	710299	98.36	ng	0.00
7) Phenol-d6	6.58	99	929006	100.57	ng	0.00
23) Nitrobenzene-d5	7.52	82	602177	70.07	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	233844	95.19	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	1073575	67.98	ng	0.00
79) Terphenyl-d14	13.07	244	880310	77.13	ng	0.00
Target Compounds						
10) Phenol	6.59	94	57506	5.824	ng	Qvalue # 64
50) Dimethylphthalate	9.70	163	174467	9.531	ng	99
52) Acenaphthene	10.03	154	32969	2.098	ng	94
71) Phenanthrene	11.51	178	170804	7.353	ng	99
72) Anthracene	11.56	178	47918	2.058	ng	98
75) Fluoranthene	12.70	202	230115	9.761	ng	98
78) Pyrene	12.93	202	166577	9.789	ng	97
81) Benzo(a)anthracene	14.12	228	39684	2.819	ng	96
83) Chrysene	14.16	228	35957	2.690	ng	98
88) Benzo(b)fluoranthene	15.20	252	31583m	2.656	ng	
90) Benzo(a)pyrene	15.59	252	22472	2.054	ng	96

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

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