

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081120\
 Data File : BF121243.D
 Acq On : 11 Aug 2020 16:46
 Operator : JU/CG
 Sample : L3625-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PN-SP

Quant Time: Aug 11 17:14:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 29 14:55:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	124183	20.00	ng	-0.01
21) Naphthalene-d8	8.07	136	498654	20.00	ng	-0.01
39) Acenaphthene-d10	9.83	164	261888	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	544656	20.00	ng	0.00
76) Chrysene-d12	13.96	240	477589	20.00	ng	0.00
86) Perylene-d12	15.42	264	424159	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	647804	85.52	ng	0.00
7) Phenol-d6	6.43	99	804997	82.27	ng	-0.02
23) Nitrobenzene-d5	7.36	82	532593	61.80	ng	-0.02
42) 2,4,6-Tribromophenol	10.63	330	236128	82.37	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1014069	57.81	ng	-0.01
79) Terphenyl-d14	12.91	244	1117988	50.89	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.55	163	75874	4.016	ng	98
87) Indeno(1,2,3-cd)pyrene	16.85	276	128591	4.359	ng	98
91) Dibenzo(a,h)anthracene	16.86	278	98026	4.068	ng	97
92) Benzo(g,h,i)perylene	17.29	276	129628	5.456	ng	99

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

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