

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062220\
 Data File : BF120695.D
 Acq On : 23 Jun 2020 00:33
 Operator : JU/CG
 Sample : L3082-16
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-17

Quant Time: Jun 23 04:13:43 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 11 11:28:22 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	143406	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	537128	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	265007	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	308330	20.00	ng	0.00
76) Chrysene-d12	13.96	240	263706	20.00	ng	0.00
86) Perylene-d12	15.40	264	393447	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	816888	97.66	ng	0.00
7) Phenol-d6	6.43	99	1009631	96.91	ng	0.00
23) Nitrobenzene-d5	7.36	82	644438	67.06	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	322979	104.83	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1123385	64.86	ng	0.00
79) Terphenyl-d14	12.91	244	951918	79.74	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.55	163	109054	5.714	ng	97
71) Phenanthrene	11.35	178	48786	3.183	ng	97
75) Fluoranthene	12.53	202	59389	3.389	ng	96
78) Pyrene	12.76	202	64532	3.454	ng	98

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

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