

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090120\
 Data File : BF121506.D
 Acq On : 1 Sep 2020 20:10
 Operator : JU/CG
 Sample : L3848-01 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LINDEN-BH-03-C

Manual Integrations
 APPROVED

mohammad
 9/2/2020 1:57:28 PM

Quant Time: Sep 02 02:16:44 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 28 09:25:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	274574	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1045443	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	514177	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	889742	20.00	ng	0.00
76) Chrysene-d12	14.03	240	678708	20.00	ng	0.00
86) Perylene-d12	15.50	264	732640	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	493433	30.15	ng	-0.01
7) Phenol-d6	6.47	99	647261	28.30	ng	-0.02
23) Nitrobenzene-d5	7.41	82	398904	26.09	ng	-0.02
42) 2,4,6-Tribromophenol	10.69	330	167736	32.30	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	836563	26.25	ng	-0.01
79) Terphenyl-d14	12.97	244	727726	22.57	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.61	163	121569	3.442	ng	99
75) Fluoranthene	12.60	202	159939	3.250	ng	95
78) Pyrene	12.83	202	162910	3.361	ng	96
81) Benzo(a)anthracene	14.02	228	88020	2.124	ng	96
88) Benzo(b)fluoranthene	15.07	252	102856m	2.158	ng	

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

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