

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062220\
 Data File : BF120689.D
 Acq On : 22 Jun 2020 21:35
 Operator : JU/CG
 Sample : L3082-18
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-19

Manual Integrations
 APPROVED

Sohil
 6/23/2020 2:03:43 PM

Quant Time: Jun 23 02:25:52 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	129869	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	513435	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	247306	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	515410	20.00	ng	0.00
76) Chrysene-d12	13.96	240	521695	20.00	ng	0.00
86) Perylene-d12	15.40	264	460201	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	786933	103.89	ng	0.00
7) Phenol-d6	6.43	99	992276	105.17	ng	0.00
23) Nitrobenzene-d5	7.36	82	606060	65.98	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	331083	115.15	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1127711	69.77	ng	0.00
79) Terphenyl-d14	12.91	244	1667954	70.62	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.55	163	136588	7.669	ng	99
75) Fluoranthene	12.53	202	117583	4.014	ng	97
78) Pyrene	12.76	202	147727	3.997	ng	99
81) Benzo(a)anthracene	13.94	228	81354	2.596	ng	98
83) Chrysene	13.98	228	74869	2.329	ng	97
88) Benzo(b)fluoranthene	14.99	252	88648m	3.234	ng	
90) Benzo(a)pyrene	15.34	252	64526	2.630	ng	98

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

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