

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011320\
 Data File : BF118541.D
 Acq On : 13 Jan 2020 20:09
 Operator : JU
 Sample : L1106-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CIY1-SB-02-03-0-32

Manual Integrations
 APPROVED

mohammad
 1/14/2020 2:31:45 PM

Quant Time: Jan 14 12:53:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 14 11:02:16 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	70895	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	248773	20.00	ng	0.00
39) Acenaphthene-d10	9.86	164	126275	20.00	ng	-0.01
64) Phenanthrene-d10	11.35	188	236961	20.00	ng	0.00
76) Chrysene-d12	14.00	240	116899	20.00	ng	-0.01
87) Perylene-d12	15.48	264	208942	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	348837	81.22	ng	0.00
7) Phenol-d6	6.46	99	369351	71.93	ng	0.00
23) Nitrobenzene-d5	7.38	82	282952	59.75	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	145606	97.72	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	450676	51.19	ng	-0.01
79) Terphenyl-d14	12.94	244	386461	52.39	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.57	163	51950	5.360	ng	99
71) Phenanthrene	11.38	178	44848	3.379	ng	98
75) Fluoranthene	12.57	202	68209	5.056	ng	99
78) Pyrene	12.80	202	67086	6.975	ng	98
80) Butylbenzylphthalate	13.41	149	12601	2.861	ng	92
81) Benzo(a)anthracene	13.99	228	23254	2.842	ng	97
83) Chrysene	14.03	228	20945	2.634	ng	99
86) Indeno(1,2,3-cd)pyrene	16.98	276	22350	3.491	ng	92
88) Benzo(b)fluoranthene	15.05	252	34150m	2.261	ng	
90) Benzo(a)pyrene	15.42	252	30299	2.448	ng	92

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

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