

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF013120\
 Data File : BF118776.D
 Acq On : 31 Jan 2020 12:39
 Operator : CG/JU
 Sample : L1318-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RAINEY-PARK-BH-04-B

Manual Integrations
 APPROVED

mohammad
 2/3/2020 4:10:14 PM

Quant Time: Jan 31 22:24:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 27 14:24:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	279810	20.00	ng	-0.04
21) Naphthalene-d8	8.11	136	1078929	20.00	ng	-0.04
39) Acenaphthene-d10	9.86	164	619163	20.00	ng	-0.04
64) Phenanthrene-d10	11.35	188	1161626	20.00	ng	-0.04
76) Chrysene-d12	13.99	240	748461	20.00	ng	-0.04
87) Perylene-d12	15.44	264	627970	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1633090	108.40	ng	-0.03
7) Phenol-d6	6.46	99	2043231	104.34	ng	-0.04
23) Nitrobenzene-d5	7.39	82	1344731	69.71	ng	-0.04
42) 2,4,6-Tribromophenol	10.65	330	813520	113.60	ng	-0.04
45) 2-Fluorobiphenyl	9.19	172	2558599	67.92	ng	-0.04
79) Terphenyl-d14	12.93	244	3000001	87.42	ng	-0.03
Target Compounds						
50) Dimethylphthalate	9.57	163	189333	4.694	ng	99
71) Phenanthrene	11.37	178	194900	3.478	ng	95
75) Fluoranthene	12.56	202	334924	5.666	ng	97
78) Pyrene	12.79	202	297652	5.858	ng	96
81) Benzo(a)anthracene	13.97	228	119593	2.656	ng	100
83) Chrysene	14.01	228	123704	2.867	ng	97
84) Bis(2-ethylhexyl)phthalate	13.97	149	77047	3.032	ng	98
88) Benzo(b)fluoranthene	15.02	252	130074m	3.295	ng	
90) Benzo(a)pyrene	15.37	252	80334	2.359	ng	# 98
92) Benzo(g,h,i)perylene	17.30	276	61116	2.090	ng	98

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

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