

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF013120\
 Data File : BF118775.D
 Acq On : 31 Jan 2020 12:11
 Operator : CG/JU
 Sample : L1318-13
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RAINEY-PARK-BH-05-A

Manual Integrations
 APPROVED

mohammad
 2/3/2020 4:10:12 PM

Quant Time: Jan 31 22:21:20 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	300750	20.00	ng	-0.04
21) Naphthalene-d8	8.11	136	1151754	20.00	ng	-0.04
39) Acenaphthene-d10	9.86	164	653097	20.00	ng	-0.04
64) Phenanthrene-d10	11.35	188	1237485	20.00	ng	-0.04
76) Chrysene-d12	13.99	240	798585	20.00	ng	-0.04
87) Perylene-d12	15.43	264	704690	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1489190	91.96	ng	-0.03
7) Phenol-d6	6.46	99	1852256	88.01	ng	-0.04
23) Nitrobenzene-d5	7.39	82	1231737	59.82	ng	-0.04
42) 2,4,6-Tribromophenol	10.65	330	708239	93.76	ng	-0.04
45) 2-Fluorobiphenyl	9.19	172	2278600	57.34	ng	-0.04
79) Terphenyl-d14	12.93	244	2611527	71.33	ng	-0.03
Target Compounds						
50) Dimethylphthalate	9.57	163	94969	2.232	ng	99
75) Fluoranthene	12.56	202	225108	3.575	ng	96
78) Pyrene	12.79	202	215521	3.975	ng	97
81) Benzo(a)anthracene	13.97	228	104428	2.174	ng	99
83) Chrysene	14.01	228	104953	2.280	ng	98
88) Benzo(b)fluoranthene	15.02	252	133112m	3.005	ng	
90) Benzo(a)pyrene	15.37	252	79478	2.080	ng	# 96

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

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