

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111819\
 Data File : BF117847.D
 Acq On : 18 Nov 2019 20:45
 Operator : JU
 Sample : K5865-11
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 3-(8-10)

Manual Integrations
 APPROVED

mohammad
 11/19/2019 2:36:28 PM

Quant Time: Nov 19 00:34:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 18:36:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	223562	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	842216	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	406684	20.00	ng	0.00
64) Phenanthrene-d10	11.46	188	659076	20.00	ng	0.00
76) Chrysene-d12	14.11	240	558048	20.00	ng	0.00
87) Perylene-d12	15.62	264	550210	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	11363	0.90	ng	0.00
7) Phenol-d6	6.53	99	317853	20.87	ng	-0.01
23) Nitrobenzene-d5	7.48	82	788884	55.68	ng	-0.01
42) 2,4,6-Tribromophenol	10.76	330	340	0.08	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	1563527	68.97	ng	0.00
79) Terphenyl-d14	13.05	244	1526107	49.98	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.67	163	162030	5.682	ng	99
71) Phenanthrene	11.49	178	271200	8.184	ng	96
75) Fluoranthene	12.68	202	306920	3.178	ng	98
78) Pyrene	12.91	202	336909	7.470	ng	98
81) Benzo(a)anthracene	14.10	228	178938	4.852	ng	96
83) Chrysene	14.14	228	175677	4.916	ng	96
88) Benzo(b)fluoranthene	15.17	252	192480m	5.384	ng	
90) Benzo(a)pyrene	15.56	252	122951	3.911	ng	# 94

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

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