

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082318\
 Data File : BF108438.D
 Acq On : 23 Aug 2018 23:16
 Operator : JU/SJ
 Sample : J4559-08 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RA-081718-SSW-145-052

Manual Integrations
 APPROVED

Sohil
 8/24/2018 12:14:20 PM

Quant Time: Aug 24 02:34:18 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 22 11:01:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	135591	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	524608	20.00	ng	0.00
38) Acenaphthene-d10	10.05	164	242940	20.00	ng	0.00
63) Phenanthrene-d10	11.54	188	384692	20.00	ng	0.00
75) Chrysene-d12	14.19	240	281577	20.00	ng	0.00
86) Perylene-d12	15.75	264	259532	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	126039	16.83	ng	-0.01
7) Phenol-d6	6.61	99	175294	17.61	ng	-0.02
23) Nitrobenzene-d5	7.56	82	110382	11.74	ng	-0.01
41) 2,4,6-Tribromophenol	10.84	330	32989	13.61	ng	-0.01
44) 2-Fluorobiphenyl	9.35	172	196416	12.98	ng	-0.01
78) Terphenyl-d14	13.12	244	128882	11.64	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.75	163	61137	3.613	ng	Qvalue # 98
70) Phenanthrene	11.57	178	220085	12.130	ng	99
71) Anthracene	11.62	178	41466	2.230	ng	96
74) Fluoranthene	12.76	202	290198	14.655	ng	95
77) Pyrene	13.00	202	237209	13.306	ng	99
80) Benzo(a)anthracene	14.18	228	109891	6.800	ng	97
82) Chrysene	14.22	228	101727	6.866	ng	98
85) Indeno(1,2,3-cd)pyrene	17.35	276	49444	3.852	ng	# 90
87) Benzo(b)fluoranthene	15.28	252	141933m	9.152	ng	
88) Benzo(k)fluoranthene	15.31	252	46657m	3.273	ng	
89) Benzo(a)pyrene	15.68	252	92361	6.651	ng	# 95
91) Benzo(a,h,i)perylene	17.85	276	46431	4.104	ng	# 87

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

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