

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082418\
 Data File : BF108482.D
 Acq On : 24 Aug 2018 23:51
 Operator : JU/SJ
 Sample : J4581-20
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP1-Q3-D4

Manual Integrations
 APPROVED

Sohil
 8/27/2018 5:02:48 PM

Quant Time: Aug 27 12:00:39 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	132635	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	485078	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	225230	20.00	ng	-0.01
63) Phenanthrene-d10	11.54	188	375320	20.00	ng	-0.01
75) Chrysene-d12	14.19	240	253669	20.00	ng	0.00
86) Perylene-d12	15.75	264	246962	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	764067	104.29	ng	0.00
7) Phenol-d6	6.62	99	1029396	105.74	ng	0.00
23) Nitrobenzene-d5	7.56	82	690647	79.45	ng	-0.01
41) 2,4,6-Tribromophenol	10.84	330	243258	108.28	ng	-0.01
44) 2-Fluorobiphenyl	9.35	172	1200015	85.54	ng	-0.01
78) Terphenyl-d14	13.12	244	889300	89.14	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.75	163	446952	28.488	ng	Qvalue # 98
70) Phenanthrene	11.56	178	72651	4.104	ng	99
74) Fluoranthene	12.76	202	114140	5.908	ng	94
77) Pyrene	12.99	202	103751	6.460	ng	98
80) Benzo(a)anthracene	14.18	228	44158	3.033	ng	# 95
82) Chrysene	14.22	228	45971m	3.444	ng	
85) Indeno(1,2,3-cd)pyrene	17.37	276	24906	2.154	ng	# 86
87) Benzo(b)fluoranthene	15.28	252	60032m	4.068	ng	
89) Benzo(a)pyrene	15.68	252	37392	2.830	ng	# 79
91) Benzo(g,h,i)perylene	17.86	276	34399	3.195	ng	# 90

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

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