

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090618\
 Data File : BF108834.D
 Acq On : 6 Sep 2018 20:15
 Operator : JU/SJ
 Sample : J4803-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP1-Q2-G5

Manual Integrations
 APPROVED

Sohil
 9/7/2018 11:15:54 AM

Quant Time: Sep 07 02:05:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF090518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 05 18:03:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	318920	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	1234086	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	614139	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	1082784	20.00	ng	0.00
75) Chrysene-d12	13.87	240	658031	20.00	ng	0.00
86) Perylene-d12	15.27	264	740049	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	1151800	59.83	ng	0.00
7) Phenol-d6	6.37	99	1542754	62.29	ng	0.00
23) Nitrobenzene-d5	7.30	82	966738	49.02	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	382201	65.19	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1741980	48.39	ng	0.00
78) Terphenyl-d14	12.82	244	1390509	49.28	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.48	163	712616	16.590	ng	99
70) Phenanthrene	11.27	178	161851	3.130	ng	98
74) Fluoranthene	12.45	202	238593	4.660	ng	97
77) Pyrene	12.67	202	228836	4.519	ng	99
80) Benzo(a)anthracene	13.85	228	99880	2.368	ng	98
82) Chrysene	13.89	228	88616	2.170	ng	97
83) Bis(2-ethylhexyl)phthalate	13.87	149	100640	3.512	ng	# 99
87) Benzo(b)fluoranthene	14.87	252	122315m	3.040	ng	
89) Benzo(a)pyrene	15.21	252	95716	2.636	ng	94

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

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