

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090718\
 Data File : BF108860.D
 Acq On : 7 Sep 2018 15:05
 Operator : JU/SJ
 Sample : J4825-17
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP1-WS-COMP-17

Manual Integrations
 APPROVED

Sohil
 9/10/2018 10:39:21 AM

Quant Time: Sep 07 15:43:04 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	214147	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	803799	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	374786	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	644324	20.00	ng	0.00
75) Chrysene-d12	13.87	240	423916	20.00	ng	0.00
86) Perylene-d12	15.28	264	538354	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1019920	78.90	ng	0.01
7) Phenol-d6	6.37	99	1317180	79.20	ng	0.00
23) Nitrobenzene-d5	7.30	82	814338	63.40	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	286399	80.05	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1448297	65.92	ng	0.00
78) Terphenyl-d14	12.82	244	1062786	58.46	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.48	163	373784	14.259	ng	99
70) Phenanthrene	11.27	178	150046	4.876	ng	98
74) Fluoranthene	12.45	202	217044	7.124	ng	98
77) Pyrene	12.68	202	209970	6.437	ng	99
80) Benzo(a)anthracene	13.86	228	103386	3.805	ng	95
82) Chrysene	13.89	228	98250m	3.735	ng	
85) Indeno(1,2,3-cd)pyrene	16.62	276	81058	3.482	ng	93
87) Benzo(b)fluoranthene	14.88	252	167991m	5.739	ng	
89) Benzo(a)pyrene	15.21	252	120967	4.580	ng	# 93
91) Benzo(g,h,i)perylene	17.04	276	78359	3.363	ng	96

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

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