

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090718\
 Data File : BF108852.D
 Acq On : 7 Sep 2018 11:31
 Operator : JU/SJ
 Sample : J4803-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP1-Q2-E7

Manual Integrations
 APPROVED

Sohil
 9/10/2018 10:39:16 AM

Quant Time: Sep 07 13:42:36 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	282317	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	1119367	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	569782	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	1089937	20.00	ng	0.00
75) Chrysene-d12	13.87	240	656114	20.00	ng	0.00
86) Perylene-d12	15.28	264	687842	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1291971	75.81	ng	0.01
7) Phenol-d6	6.37	99	1705003	77.76	ng	0.00
23) Nitrobenzene-d5	7.30	82	1073802	60.03	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	424103	77.97	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1827864	54.73	ng	0.00
78) Terphenyl-d14	12.83	244	1716296	61.00	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.48	163	810512	20.338	ng	99
70) Phenanthrene	11.27	178	110994	2.132	ng	98
74) Fluoranthene	12.45	202	238663	4.631	ng	99
77) Pyrene	12.67	202	224239	4.442	ng	98
82) Chrysene	13.89	228	88609	2.176	ng	97
83) Bis(2-ethylhexyl)phthalate	13.87	149	74065	2.592	ng	99
87) Benzo(b)fluoranthene	14.88	252	115291m	3.083	ng	
89) Benzo(a)pyrene	15.21	252	76181	2.257	ng	# 91
91) Benzo(g,h,i)perylene	17.03	276	61524	2.066	ng	96

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

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