

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090518\
 Data File : BF108810.D
 Acq On : 6 Sep 2018 8:04
 Operator : JU/SJ
 Sample : J4724-04
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP1-SUT-UST-03

Manual Integrations
 APPROVED

Sohil
 9/6/2018 4:46:12 PM

Quant Time: Sep 06 11:14:46 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	283372	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	1067584	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	501140	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	834646	20.00	ng	0.00
75) Chrysene-d12	13.87	240	610817	20.00	ng	0.00
86) Perylene-d12	15.27	264	596644	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	977528	57.15	ng	0.00
7) Phenol-d6	6.37	99	1262108	57.35	ng	0.00
23) Nitrobenzene-d5	7.29	82	768792	45.07	ng	-0.01
41) 2,4,6-Tribromophenol	10.55	330	273450	57.16	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1423591	48.46	ng	0.00
78) Terphenyl-d14	12.82	244	1096712	41.87	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.48	163	553735	15.798	ng	99
70) Phenanthrene	11.27	178	284579	7.140	ng	97
74) Fluoranthene	12.45	202	382563	9.694	ng	97
77) Pyrene	12.67	202	341898	7.274	ng	99
80) Benzo(a)anthracene	13.85	228	171737	4.386	ng	98
82) Chrysene	13.89	228	151799	4.005	ng	95
85) Indeno(1,2,3-cd)pyrene	16.62	276	76283	2.274	ng	95
87) Benzo(b)fluoranthene	14.87	252	190919m	5.885	ng	
88) Benzo(k)fluoranthene	14.89	252	63969m	2.141	ng	
89) Benzo(a)pyrene	15.21	252	132179	4.515	ng	98
91) Benzo(g,h,i)perylene	17.02	276	71438	2.766	ng	98

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

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