

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111618\
 Data File : BF110815.D
 Acq On : 16 Nov 2018 14:33
 Operator : JU/SJ
 Sample : J5964-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LR-RBR-200029

Manual Integrations
 APPROVED

Sohil
 11/19/2018 3:48:10 PM

Quant Time: Nov 16 15:39:51 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	63589	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	265613	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	127455	20.00	ng	-0.01
64) Phenanthrene-d10	11.47	188	236931	20.00	ng	-0.01
76) Chrysene-d12	14.13	240	140861	20.00	ng	-0.01
87) Perylene-d12	15.66	264	144611	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	487541	124.54	ng	0.00
7) Phenol-d6	6.57	99	624505	124.75	ng	0.00
23) Nitrobenzene-d5	7.50	82	394400	85.51	ng	-0.01
42) 2,4,6-Tribromophenol	10.78	330	142581	111.02	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	737503	82.64	ng	0.00
79) Terphenyl-d14	13.06	244	596337	79.28	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.69	163	73053	7.681	ng	99
71) Phenanthrene	11.50	178	33770	2.660	ng	98
75) Fluoranthene	12.70	202	54644	4.294	ng	98
78) Pyrene	12.93	202	58541	5.397	ng	98
81) Benzo(a)anthracene	14.12	228	22930	2.723	ng	98
83) Chrysene	14.16	228	22845	2.848	ng	99
88) Benzo(b)fluoranthene	15.20	252	20465m	2.366	ng	
90) Benzo(a)pyrene	15.59	252	16345	2.079	ng	99

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

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