

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
 Data File : BF110530.D
 Acq On : 6 Nov 2018 22:07
 Operator : JU/SJ
 Sample : J5764-05 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-03-110218-C

Manual Integrations
 APPROVED

Sohil
 11/7/2018 2:54:42 PM

Quant Time: Nov 07 12:43:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 16:06:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	86786	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	329927	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	125354	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	198547	20.00	ng	0.00
76) Chrysene-d12	14.14	240	154645	20.00	ng	0.00
87) Perylene-d12	15.66	264	107560	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	254799	47.81	ng	0.00
7) Phenol-d6	6.57	99	330204	48.44	ng	0.00
23) Nitrobenzene-d5	7.52	82	188765	33.94	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	51487	41.46	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	356539	44.66	ng	0.00
79) Terphenyl-d14	13.07	244	273760	36.82	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.70	163	44187	4.775	ng	Qvalue 100
71) Phenanthrene	11.52	178	24674	2.375	ng	98
75) Fluoranthene	12.71	202	41709	3.956	ng	99
78) Pyrene	12.94	202	39120	3.529	ng	98
81) Benzo(a)anthracene	14.13	228	20796	2.268	ng	93
83) Chrysene	14.16	228	18333	2.105	ng	95
88) Benzo(b)fluoranthene	15.21	252	15883m	2.557	ng	
90) Benzo(a)pyrene	15.60	252	14006	2.451	ng	# 77

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

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