

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091018\
 Data File : BF108907.D
 Acq On : 10 Sep 2018 21:38
 Operator : JU/SJ
 Sample : J4821-06
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ENY-BC-02

Manual Integrations
 APPROVED

Sohil
 9/11/2018 9:34:17 AM

Quant Time: Sep 11 03:59:29 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	196036	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	707222	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	304338	20.00	ng	0.00
63) Phenanthrene-d10	11.25	188	495029	20.00	ng	0.00
75) Chrysene-d12	13.88	240	410829	20.00	ng	0.01
86) Perylene-d12	15.29	264	268820	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1110537	93.85	ng	0.02
7) Phenol-d6	6.38	99	1432793	94.11	ng	0.01
23) Nitrobenzene-d5	7.30	82	930138	82.31	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	263169	90.58	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1457789	81.72	ng	0.00
78) Terphenyl-d14	12.83	244	1134957	64.42	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.49	163	451646	21.218	ng	99
74) Fluoranthene	12.45	202	62310	2.662	ng	96
77) Pyrene	12.68	202	67264	2.128	ng	99
87) Benzo(b)fluoranthene	14.89	252	47204m	3.229	ng	
89) Benzo(a)pyrene	15.23	252	30007	2.275	ng	# 73
91) Benzo(g,h,i)perylene	17.06	276	27493	2.363	ng	# 92

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

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