

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100518\
 Data File : BF109622.D
 Acq On : 5 Oct 2018 21:32
 Operator : JU/SJ
 Sample : J5255-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRACK-D-2

Manual Integrations
 APPROVED

Sohil
 10/8/2018 12:21:27 PM

Quant Time: Oct 06 06:45:02 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	83032	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	321170	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	142747	20.00	ng	-0.01
63) Phenanthrene-d10	11.21	188	229807	20.00	ng	0.00
75) Chrysene-d12	13.83	240	193952	20.00	ng	-0.01
86) Perylene-d12	15.23	264	200922	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	465556	92.35	ng	0.01
7) Phenol-d6	6.35	99	617495	95.99	ng	0.00
23) Nitrobenzene-d5	7.26	82	401689	73.83	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	127455	90.57	ng	-0.02
44) 2-Fluorobiphenyl	9.06	172	727818	76.90	ng	-0.02
78) Terphenyl-d14	12.79	244	525725	66.52	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.45	163	67708	6.521	ng	99
74) Fluoranthene	12.42	202	38520	3.141	ng	95
77) Pyrene	12.64	202	37307	2.684	ng	99
80) Benzo(a)anthracene	13.82	228	24231	2.074	ng	94
82) Chrysene	13.86	228	27673	2.390	ng	99
87) Benzo(b)fluoranthene	14.84	252	43286m	3.921	ng	
89) Benzo(a)pyrene	15.18	252	20054	2.016	ng	95

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

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