

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
 Data File : BF107342.D
 Acq On : 12 Jul 2018 16:42
 Operator : JU/SJ
 Sample : J3930-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NYS-RBR-200030

Manual Integrations
 APPROVED

Sohil
 7/13/2018 12:37:44 PM

Quant Time: Jul 13 08:33:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 11 16:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	74002	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	285365	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	139861	20.00	ng	0.00
63) Phenanthrene-d10	11.48	188	247319	20.00	ng	0.00
75) Chrysene-d12	14.14	240	212089	20.00	ng	0.00
86) Perylene-d12	15.68	264	230006	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	404555	92.57	ng	0.00
7) Phenol-d6	6.58	99	511502	93.72	ng	0.00
23) Nitrobenzene-d5	7.50	82	327911	63.86	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	149002	91.92	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	580911	67.22	ng	0.00
78) Terphenyl-d14	13.06	244	565517	64.59	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.68	163	128788	12.278	ng	98
70) Phenanthrene	11.49	178	28316	2.374	ng	98
74) Fluoranthene	12.69	202	62128	4.725	ng	96
77) Pyrene	12.93	202	56162	4.016	ng	99
80) Benzo(a)anthracene	14.12	228	33618	2.601	ng	100
82) Chrysene	14.16	228	36854	3.099	ng	99
85) Indeno(1,2,3-cd)pyrene	17.27	276	21156m	2.096	ng	
87) Benzo(b)fluoranthene	15.22	252	53883m	3.771	ng	
89) Benzo(a)pyrene	15.61	252	34809	2.741	ng	98

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

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