

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF040218\
 Data File : BF104149.D
 Acq On : 2 Apr 2018 21:22
 Operator : JU/SJ
 Sample : J2112-02
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-1-COMP

Manual Integrations
 APPROVED

Sohil
 4/3/2018 2:27:38 PM

Quant Time: Apr 03 02:09:22 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 02 13:40:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	96354	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	356630	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	142623	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	218103	20.00	ng	0.00
75) Chrysene-d12	14.15	240	237604	20.00	ng	0.00
86) Perylene-d12	15.70	264	201057	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	683392	113.11	ng	0.00
7) Phenol-d6	6.59	99	808890	108.82	ng	0.00
23) Nitrobenzene-d5	7.52	82	502216	86.12	ng	0.00
41) 2,4,6-Tribromophenol	10.79	330	161871	101.93	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	754052	83.37	ng	0.00
78) Terphenyl-d14	13.08	244	634926	61.70	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.70	163	57209	5.080	ng	Qvalue 100
70) Phenanthrene	11.52	178	41126	3.483	ng	98
74) Fluoranthene	12.72	202	70593	5.624	ng	96
77) Pyrene	12.95	202	66331	3.883	ng	95
80) Benzo(a)anthracene	14.14	228	42141	2.834	ng	95
82) Chrysene	14.17	228	34957	2.401	ng	97
87) Benzo(b)fluoranthene	15.23	252	40321m	3.295	ng	
89) Benzo(a)pyrene	15.63	252	29086	2.565	ng	# 78

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

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