

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073118\
 Data File : BF107864.D
 Acq On : 31 Jul 2018 21:54
 Operator : JU/SJ
 Sample : J4236-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-01-010-A

Manual Integrations
 APPROVED

Sohil
 8/1/2018 3:57:28 PM

Quant Time: Aug 01 03:31:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 17:48:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	157991	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	664137	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	287321	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	515775	20.00	ng	0.00
75) Chrysene-d12	14.15	240	438140	20.00	ng	0.00
86) Perylene-d12	15.68	264	443915	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	973569	104.73	ng	0.00
7) Phenol-d6	6.60	99	1194116	97.47	ng	0.00
23) Nitrobenzene-d5	7.52	82	802015	69.57	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	320597	82.12	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	1492915	72.41	ng	0.00
78) Terphenyl-d14	13.08	244	1169470	57.88	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	99769	4.624	ng	99
70) Phenanthrene	11.52	178	70101	2.862	ng	96
74) Fluoranthene	12.71	202	130643	4.541	ng	98
77) Pyrene	12.95	202	127016	4.026	ng	97
80) Benzo(a)anthracene	14.14	228	67483	2.719	ng	98
82) Chrysene	14.18	228	62350	2.329	ng	97
87) Benzo(b)fluoranthene	15.23	252	73675m	3.365	ng	
89) Benzo(a)pyrene	15.62	252	57924	2.563	ng	# 87

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

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