

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
 Data File : BF107643.D
 Acq On : 25 Jul 2018 3:58
 Operator : JU/SJ
 Sample : J4120-04
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ST-3

Manual Integrations
 APPROVED

Sohil
 7/25/2018 12:58:46 PM

Quant Time: Jul 25 08:32:16 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 16:29:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	256803	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	959755	20.00	ng	-0.01
38) Acenaphthene-d10	10.01	164	367171	20.00	ng	-0.01
63) Phenanthrene-d10	11.50	188	629947	20.00	ng	-0.01
75) Chrysene-d12	14.15	240	486674	20.00	ng	-0.01
86) Perylene-d12	15.68	264	377271	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	1421707	89.01	ng	0.00
7) Phenol-d6	6.60	99	1728651	90.13	ng	-0.01
23) Nitrobenzene-d5	7.53	82	1050993	73.90	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	354060	84.80	ng	-0.02
44) 2-Fluorobiphenyl	9.32	172	1698600	77.26	ng	-0.01
78) Terphenyl-d14	13.09	244	1387670	73.00	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.71	163	100758	3.641	ng	99
70) Phenanthrene	11.52	178	156680	5.104	ng	98
74) Fluoranthene	12.72	202	160878	4.884	ng	94
77) Pyrene	12.95	202	202871	6.094	ng	96
80) Benzo(a)anthracene	14.14	228	77311m	2.711	ng	
82) Chrysene	14.18	228	77619	2.833	ng	95
87) Benzo(b)fluoranthene	15.23	252	48465m	2.347	ng	
89) Benzo(a)pyrene	15.62	252	41552	2.200	ng	94

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

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