

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022718\
 Data File : BF103260.D
 Acq On : 27 Feb 2018 14:02
 Operator : SJ/JU
 Sample : J1702-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-03-003-A

Manual Integrations
 APPROVED

Sohil
 2/28/2018 11:17:30 AM

Quant Time: Feb 27 17:13:09 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 23 02:43:46 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	158722	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	665399	20.00	ng	0.00
38) Acenaphthene-d10	9.91	164	293303	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	476129	20.00	ng	0.00
75) Chrysene-d12	14.05	240	311896	20.00	ng	0.00
86) Perylene-d12	15.55	264	328399	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1009119	96.16	ng	0.01
7) Phenol-d6	6.51	99	1214045	94.54	ng	0.00
23) Nitrobenzene-d5	7.44	82	793278	68.08	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	235056	94.68	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1073421	58.39	ng	0.00
78) Terphenyl-d14	12.99	244	817862	63.03	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	149416	6.350	ng	98
70) Phenanthrene	11.42	178	63596	2.427	ng	97
74) Fluoranthene	12.62	202	110340	4.191	ng	98
77) Pyrene	12.84	202	110486	4.544	ng	97
80) Benzo(a)anthracene	14.04	228	45498	2.248	ng	99
82) Chrysene	14.07	228	44737	2.277	ng	98
87) Benzo(b)fluoranthene	15.11	252	63133m	2.924	ng	
89) Benzo(a)pyrene	15.48	252	44828	2.325	ng	# 89

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

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