

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080118\
 Data File : BF107886.D
 Acq On : 1 Aug 2018 13:59
 Operator : JU/SJ
 Sample : J4237-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-023-A

Manual Integrations
 APPROVED

Sohil
 8/2/2018 9:17:00 AM

Quant Time: Aug 01 14:56:58 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	6.97	152	132862	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	540749	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	232019	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	426724	20.00	ng	0.00
75) Chrysene-d12	14.15	240	400971	20.00	ng	0.00
86) Perylene-d12	15.69	264	451593	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	785727	100.51	ng	0.00
7) Phenol-d6	6.60	99	908670	88.19	ng	0.00
23) Nitrobenzene-d5	7.52	82	616465	65.68	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	245074	77.74	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	1070699	64.31	ng	0.00
78) Terphenyl-d14	13.08	244	952319	51.50	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.71	163	179354	10.293	ng	Qvalue 100
70) Phenanthrene	11.52	178	187302	9.244	ng	98
71) Anthracene	11.58	178	74421	3.184	ng	96
74) Fluoranthene	12.72	202	197045	8.278	ng	96
77) Pyrene	12.95	202	168154	5.824	ng	95
80) Benzo(a)anthracene	14.14	228	90214	3.972	ng	97
82) Chrysene	14.18	228	74740	3.051	ng	97
85) Indeno(1,2,3-cd)pyrene	17.28	276	41241	2.213	ng	# 90
87) Benzo(b)fluoranthene	15.24	252	89830m	4.033	ng	
89) Benzo(a)pyrene	15.63	252	72373	3.148	ng	# 94

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

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