

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
 Data File : BF127192.D
 Acq On : 04 Mar 2022 01:08
 Operator : CG\JU
 Sample : N1725-03DL 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-03DL

Quant Time: Mar 04 02:16:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 01 23:57:09 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.886	152	101186	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	386134	20.000	ng	# 0.00
39) Acenaphthene-d10	9.933	164	188096	20.000	ng	0.00
64) Phenanthrene-d10	11.421	188	278400	20.000	ng	# 0.00
76) Chrysene-d12	14.068	240	192138	20.000	ng	# 0.00
86) Perylene-d12	15.562	264	159355	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	49196	7.514	ng	0.00
7) Phenol-d6	6.516	99	43110	4.880	ng	0.00
23) Nitrobenzene-d5	7.451	82	135050	15.787	ng	-0.01
42) 2,4,6-Tribromophenol	10.722	330	26885	12.414	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	213222	16.688	ng	0.00
79) Terphenyl-d14	13.004	244	177424	13.575	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.198	128	662098	32.847	ng	99
37) 2-Methylnaphthalene	8.886	142	44563	3.407	ng	95
38) 1-Methylnaphthalene	8.986	142	30884	2.466	ng	92
52) Acenaphthene	9.963	154	65642	5.623	ng	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030322\
Data File : BF127192.D
Acq On : 04 Mar 2022 01:08
Operator : CG\JU
Sample : N1725-03DL 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-03DL

Quant Time: Mar 04 02:16:05 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 01 23:57:09 2022
Response via : Initial Calibration

