

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073021\
 Data File : BF124880.D
 Acq On : 30 Jul 2021 17:19
 Operator : JU/CG
 Sample : M3229-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SOLID-A-E

Quant Time: Jul 30 17:42:57 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 22 15:28:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	6269	20.000	ng	# 0.00
21) Naphthalene-d8	8.157	136	26493	20.000	ng	# 0.00
39) Acenaphthene-d10	9.916	164	16093	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	26927	20.000	ng	# 0.00
76) Chrysene-d12	14.039	240	14379	20.000	ng	# 0.00
86) Perylene-d12	15.515	264	12982	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	34130	92.180	ng	0.00
7) Phenol-d6	6.498	99	42369	91.279	ng	0.00
23) Nitrobenzene-d5	7.434	82	29497	66.266	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	12624	91.363	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	60681	55.381	ng	0.00
79) Terphenyl-d14	12.986	244	59942	71.458	ng	0.00
Target Compounds						
						Qvalue
37) 2-Methylnaphthalene	8.869	142	5745	6.219	ng	92
38) 1-Methylnaphthalene	8.969	142	24855	28.401	ng	100
58) Fluorene	10.469	166	6263	5.877	ng	# 90
71) Phenanthrene	11.427	178	15755	10.879	ng	# 95

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

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