

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062922\
 Data File : BG054164.D
 Acq On : 30 Jun 2022 3:29
 Operator : CG/JU
 Sample : N3527-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 600532979-B

Quant Time: Jun 30 05:58:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 27 00:27:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.065	152	22483	20.000	ng	0.00
21) Naphthalene-d8	10.879	136	89633	20.000	ng	# 0.00
39) Acenaphthene-d10	14.693	164	65927	20.000	ng	0.00
64) Phenanthrene-d10	17.436	188	156966	20.000	ng	# 0.00
76) Chrysene-d12	21.714	240	172111	20.000	ng	0.00
86) Perylene-d12	24.975	264	179447	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.638	112	52761	44.867	ng	0.00
7) Phenol-d6	7.231	99	113708	71.279	ng	0.00
23) Nitrobenzene-d5	9.228	82	110187	69.083	ng	0.00
42) 2,4,6-Tribromophenol	16.179	330	26748	25.252	ng	0.00
45) 2-Fluorobiphenyl	13.318	172	320108	70.605	ng	0.00
79) Terphenyl-d14	20.045	244	609684	71.347	ng	0.00
Target Compounds						
37) 2-Methylnaphthalene	12.530	142	11992	3.582	ng	96
38) 1-Methylnaphthalene	12.748	142	8121	2.491	ng	98

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

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