

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042924\
 Data File : BG061051.D
 Acq On : 29 Apr 2024 17:39
 Operator : MA/JU
 Sample : P2270-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 72-11936

Quant Time: Apr 29 18:35:07 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 22 04:22:00 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.943	152	21801	20.000	ng	0.00
21) Naphthalene-d8	10.734	136	82958	20.000	ng	# 0.00
39) Acenaphthene-d10	14.570	164	74214	20.000	ng	0.00
64) Phenanthrene-d10	17.314	188	200051	20.000	ng	0.00
76) Chrysene-d12	21.574	240	206226	20.000	ng	# 0.00
86) Perylene-d12	24.694	264	229710	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	160083	148.529	ng	0.00
7) Phenol-d6	7.114	99	195235	122.087	ng	0.00
23) Nitrobenzene-d5	9.094	82	165026	89.376	ng	0.00
42) 2,4,6-Tribromophenol	16.063	330	237544	145.761	ng	0.00
45) 2-Fluorobiphenyl	13.195	172	496063	92.510	ng	0.00
79) Terphenyl-d14	19.940	244	1313323	98.852	ng	0.00

Target Compounds	Qvalue

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

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