

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122223\
 Data File : BG060161.D
 Acq On : 23 Dec 2023 7:01
 Operator : MA/JU
 Sample : 05862-01
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-LIQUIDS-20231213-01

Quant Time: Dec 25 00:44:21 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 03:42:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.941	152	318844	20.000	ng	0.00
21) Naphthalene-d8	10.744	136	1431095	20.000	ng	0.00
39) Acenaphthene-d10	14.575	164	1070032	20.000	ng	0.00
64) Phenanthrene-d10	17.325	188	2316261	20.000	ng	0.00
76) Chrysene-d12	21.584	240	1583815	20.000	ng	-0.01
86) Perylene-d12	24.763	264	1789767	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.515	112	2341644	115.723	ng	0.00
7) Phenol-d6	7.107	99	3331148	97.818	ng	0.00
23) Nitrobenzene-d5	9.105	82	2296729	78.619	ng	0.00
42) 2,4,6-Tribromophenol	16.067	330	1482576	102.906	ng	0.00
45) 2-Fluorobiphenyl	13.200	172	5084777	71.353	ng	0.00
79) Terphenyl-d14	19.945	244	5348928	61.112	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.803	128	7201729	94.530	ng	# 79
37) 2-Methylnaphthalene	12.401	142	1067286	18.051	ng	99
38) 1-Methylnaphthalene	12.624	142	819928	13.815	ng	98
52) Acenaphthene	14.640	154	145513	2.374	ng	92
71) Phenanthrene	17.366	178	255287	2.187	ng	95

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

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