

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG103122\
 Data File : BG055386.D
 Acq On : 31 Oct 2022 19:53
 Operator : CG/JU
 Sample : N5377-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LIQUID-WC-20221028

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/01/2022
 Supervised By : mohammad ahmed 11/04/2022

Quant Time: Nov 01 06:50:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 29 04:52:38 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.259	152	33890	20.000	ng	0.00
21) Naphthalene-d8	11.091	136	137228	20.000	ng	# 0.00
39) Acenaphthene-d10	14.874	164	93722	20.000	ng	0.00
64) Phenanthrene-d10	17.606	188	201745	20.000	ng	# 0.00
76) Chrysene-d12	21.884	240	194787	20.000	ng	0.00
86) Perylene-d12	25.274	264	207843	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.791	112	182073	94.070	ng	0.00
7) Phenol-d6	7.407	99	279014	99.503	ng	0.00
23) Nitrobenzene-d5	9.434	82	217824	81.054	ng	0.00
42) 2,4,6-Tribromophenol	16.379	330	2606m	2.558	ng	0.02
45) 2-Fluorobiphenyl	13.500	172	522353	91.223	ng	-0.01
79) Terphenyl-d14	20.180	244	745651	79.535	ng	0.00

Target Compounds Qvalue

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

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