

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052822\
 Data File : BG053753.D
 Acq On : 28 May 2022 16:31
 Operator : CG/JU
 Sample : N3056-09DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-02DL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/29/2022
 Supervised By : Jagrut Upadhyay 05/31/2022

Quant Time: May 29 02:19:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun May 29 02:17:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.947	152	37826	20.000	ng	0.00
21) Naphthalene-d8	10.755	136	153726	20.000	ng	0.00
39) Acenaphthene-d10	14.591	164	110244	20.000	ng	0.00
64) Phenanthrene-d10	17.340	188	271877	20.000	ng	0.00
76) Chrysene-d12	21.599	240	285682	20.000	ng	-0.01
86) Perylene-d12	24.777	264	253710	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.526	112	27447	12.315	ng	0.00
7) Phenol-d6	7.154	99	26678	7.887	ng	0.00
23) Nitrobenzene-d5	9.116	82	60857	16.852	ng	0.00
42) 2,4,6-Tribromophenol	16.089	330	43025	26.461	ng	0.00
45) 2-Fluorobiphenyl	13.210	172	149678	19.840	ng	0.00
79) Terphenyl-d14	19.948	244	220767	14.393	ng	0.00
Target Compounds						Qvalue
27) 2,4-Dimethylphenol	9.956	122	5328m	2.501	ng	
31) Naphthalene	10.807	128	350842	44.867	ng	99
37) 2-Methylnaphthalene	12.417	142	50649	8.679	ng	97
38) 1-Methylnaphthalene	12.634	142	45330m	7.962	ng	

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

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