

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020824\
 Data File : BF137353.D
 Acq On : 08 Feb 2024 22:44
 Operator : CG\JU
 Sample : P1372-05DL2 100X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CG-MW-17-ch4DL2

Quant Time: Feb 09 02:23:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF013024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 31 04:56:22 2024
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 02/09/2024
 Supervised By :mohammad ahmed 02/09/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.787	152	84908	20.000	ng	0.00
21) Naphthalene-d8	8.063	136	323973	20.000	ng	#-0.01
39) Acenaphthene-d10	9.816	164	169006	20.000	ng	-0.01
64) Phenanthrene-d10	11.310	188	272743	20.000	ng	0.00
76) Chrysene-d12	13.963	240	188957	20.000	ng	-0.01
86) Perylene-d12	15.445	264	209898	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	3095	0.601	ng	0.02
7) Phenol-d6	6.440	99	5116m	0.697	ng	0.01
23) Nitrobenzene-d5	7.351	82	14238	1.968	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	2009	1.053	ng	0.00
45) 2-Fluorobiphenyl	9.139	172	37272	3.032	ng	-0.01
79) Terphenyl-d14	12.910	244	33073	2.379	ng	-0.01
Target Compounds						
31) Naphthalene	8.092	128	2030813	115.951	ng	99
38) 1-Methylnaphthalene	8.875	142	145980	13.385	ng	98
52) Acenaphthene	9.851	154	32287	3.448	ng	98

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

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