

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010522\
 Data File : BP008694.D
 Acq On : 05 Jan 2022 17:27
 Operator : CG/JU
 Sample : M5251-01DL 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-1RDL

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 01/06/2022
 Supervised By :Jagrut Upadhyay 01/06/2022

Quant Time: Jan 06 03:03:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP122821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 30 08:44:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	560263	20.000	ng	-0.02
21) Naphthalene-d8	10.687	136	2241651	20.000	ng	-0.01
39) Acenaphthene-d10	14.516	164	1420066	20.000	ng	-0.01
64) Phenanthrene-d10	17.263	188	2942129	20.000	ng	-0.01
76) Chrysene-d12	21.351	240	2614706	20.000	ng	-0.01
86) Perylene-d12	23.774	264	2798225	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	162374	5.290	ng	0.00
7) Phenol-d6	7.052	99	155208	3.633	ng	0.00
23) Nitrobenzene-d5	9.040	82	260937	6.925	ng	-0.02
42) 2,4,6-Tribromophenol	16.004	330	194429	10.120	ng	-0.01
45) 2-Fluorobiphenyl	13.140	172	777935	7.937	ng	-0.02
79) Terphenyl-d14	19.822	244	849306	6.632	ng	0.00
Target Compounds						Qvalue
27) 2,4-Dimethylphenol	9.893	122	117885m	3.833	ng	
31) Naphthalene	10.734	128	5409303	46.491	ng	99
37) 2-Methylnaphthalene	12.345	142	644528	7.359	ng	99
38) 1-Methylnaphthalene	12.563	142	348074m	4.065	ng	

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

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