

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031824\
 Data File : BF137630.D
 Acq On : 18 Mar 2024 14:37
 Operator : CG\JU
 Sample : P1769-01DL2 250X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2DL2

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 03/19/2024
 Supervised By :mohammad ahmed 03/19/2024

Quant Time: Mar 18 15:09:03 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 14 14:45:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.728	152	116165	20.000	ng	0.00
21) Naphthalene-d8	8.004	136	455084	20.000	ng	0.00
39) Acenaphthene-d10	9.757	164	261489	20.000	ng	0.00
64) Phenanthrene-d10	11.239	188	464790	20.000	ng	0.00
76) Chrysene-d12	13.874	240	314801	20.000	ng	0.00
86) Perylene-d12	15.298	264	269931	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	0.000	112	0d	0.000	ng	
7) Phenol-d6	6.363	99	1572	0.142	ng	0.00
23) Nitrobenzene-d5	7.304	82	3412m	0.348	ng	0.00
42) 2,4,6-Tribromophenol	10.563	330	1105	0.481	ng	0.01
45) 2-Fluorobiphenyl	9.086	172	9246	0.553	ng	0.00
79) Terphenyl-d14	12.827	244	11384	0.497	ng	0.00
Target Compounds						
10) Phenol	6.369	94	324017	27.191	ng	Qvalue 93
17) 2-Methylphenol	6.981	107	91013	12.023	ng	93
20) 3+4-Methylphenols	7.139	107	75979	8.764	ng	# 53
27) 2,4-Dimethylphenol	7.669	122	36599	5.014	ng	96

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

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