

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022723\
 Data File : BF132547.D
 Acq On : 27 Feb 2023 12:14
 Operator : CG\JU
 Sample : 01654-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 UGT

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 02/28/2023
 Supervised By :Jagrut Upadhyay 02/28/2023

Quant Time: Feb 27 13:19:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 24 23:17:30 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.001	152	149452	20.000	ng	0.00
21) Naphthalene-d8	8.289	136	540219	20.000	ng	0.00
39) Acenaphthene-d10	10.048	164	298629	20.000	ng	0.00
64) Phenanthrene-d10	11.542	188	575648	20.000	ng	0.00
76) Chrysene-d12	14.201	240	314970	20.000	ng	0.00
86) Perylene-d12	15.754	264	306605	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.643	112	1158918	125.852	ng	0.01
7) Phenol-d6	6.631	99	1395183	116.091	ng	0.00
23) Nitrobenzene-d5	7.566	82	1095645	96.224	ng	0.00
42) 2,4,6-Tribromophenol	10.848	330	475584	147.465	ng	0.00
45) 2-Fluorobiphenyl	9.366	172	1743501	88.898	ng	0.00
79) Terphenyl-d14	13.136	244	2017408	108.294	ng	0.00
Target Compounds						
32) Benzoic acid	8.037	122	56770m	11.652	ng	Qvalue

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

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