

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022522\
 Data File : BM034087.D
 Acq On : 25 Feb 2022 20:50
 Operator : CG/JU
 Sample : N1664-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-3

Quant Time: Feb 26 04:02:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 26 03:45:21 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 02/28/2022
 Supervised By :Jagrut Upadhyay 02/28/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.001	152	124926	20.000	ng	0.00
21) Naphthalene-d8	10.819	136	441953	20.000	ng	# 0.00
39) Acenaphthene-d10	14.636	164	276484	20.000	ng	0.00
64) Phenanthrene-d10	17.377	188	565951	20.000	ng	# 0.00
76) Chrysene-d12	21.547	240	534686	20.000	ng	# 0.00
86) Perylene-d12	24.000	264	530173	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.554	112	992421	141.613	ng	0.00
7) Phenol-d6	7.154	99	1064811	115.709	ng	0.00
23) Nitrobenzene-d5	9.166	82	819788	100.101	ng	0.00
42) 2,4,6-Tribromophenol	16.118	330	600797	179.042	ng	0.00
45) 2-Fluorobiphenyl	13.266	172	2181419	105.116	ng	0.00
79) Terphenyl-d14	19.995	244	3398912	110.694	ng	0.00
Target Compounds						
32) Benzoic acid	10.036	122	14008m	3.187	ng	Qvalue

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

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