

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080423\
 Data File : BM041296.D
 Acq On : 04 Aug 2023 16:09
 Operator : MA/JU
 Sample : 03887-08
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MCLEAN-TP4

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 08/07/2023
 Supervised By :mohammad ahmed 08/07/2023

Quant Time: Aug 04 17:49:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM072823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 02 19:20:51 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.792	152	102399	20.000	ng	0.00
21) Naphthalene-d8	10.598	136	427073	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	300879	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	822959	20.000	ng	0.00
76) Chrysene-d12	21.421	240	849395	20.000	ng	-0.01
86) Perylene-d12	23.797	264	833944	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	786322	114.769	ng	0.00
7) Phenol-d6	6.975	99	981654	110.535	ng	0.00
23) Nitrobenzene-d5	8.969	82	716286	77.997	ng	0.00
42) 2,4,6-Tribromophenol	15.962	330	672219	157.955	ng	0.00
45) 2-Fluorobiphenyl	13.074	172	1692200	71.047	ng	0.00
79) Terphenyl-d14	19.856	244	4387943	93.403	ng	0.00
Target Compounds						
32) Benzoic acid	9.951	122	9193m	8.786	ng	Qvalue

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

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