

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080423\
 Data File : BM041295.D
 Acq On : 04 Aug 2023 15:32
 Operator : MA/JU
 Sample : 03887-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MCLEAN-TP7

Manual Integrations
 APPROVED

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

Quant Time: Aug 04 16:14:13 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	136001	20.000	ng	0.00
21) Naphthalene-d8	10.598	136	559986	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	376565	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	926592	20.000	ng	0.00
76) Chrysene-d12	21.421	240	833859	20.000	ng	-0.01
86) Perylene-d12	23.797	264	809984	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	1069990	117.587	ng	0.00
7) Phenol-d6	6.975	99	1298838	110.115	ng	0.00
23) Nitrobenzene-d5	8.969	82	963740	80.035	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	786970	147.751	ng	0.00
45) 2-Fluorobiphenyl	13.074	172	2164065	72.597	ng	0.00
79) Terphenyl-d14	19.856	244	4596784	99.672	ng	0.00
Target Compounds						
32) Benzoic acid	9.957	122	12753m	8.880	ng	Qvalue

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

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