

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080423\
 Data File : BM041294.D
 Acq On : 04 Aug 2023 14:55
 Operator : MA/JU
 Sample : 03887-14
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MCLEAN-TP9

Manual Integrations
 APPROVED

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

Quant Time: Aug 04 16:14:09 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	129530	20.000	ng	0.00
21) Naphthalene-d8	10.598	136	536546	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	376979	20.000	ng	0.00
64) Phenanthrene-d10	17.215	188	999726	20.000	ng	0.00
76) Chrysene-d12	21.421	240	982874	20.000	ng	-0.01
86) Perylene-d12	23.797	264	949178	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	983585	113.491	ng	0.00
7) Phenol-d6	6.975	99	1181572	105.178	ng	0.00
23) Nitrobenzene-d5	8.969	82	927016	80.348	ng	0.00
42) 2,4,6-Tribromophenol	15.962	330	836127	156.808	ng	0.00
45) 2-Fluorobiphenyl	13.074	172	2197365	73.633	ng	0.00
79) Terphenyl-d14	19.856	244	5253864	96.648	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	9.951	122	14588m	9.215	ng	

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

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