

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080423\
 Data File : BM041290.D
 Acq On : 04 Aug 2023 12:28
 Operator : MA/JU
 Sample : 03877-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COMP-2

Manual Integrations
 APPROVED

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

Quant Time: Aug 04 16:13:53 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	122374	20.000	ng	0.00
21) Naphthalene-d8	10.598	136	504873	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	354168	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	925326	20.000	ng	# 0.00
76) Chrysene-d12	21.421	240	863521	20.000	ng	-0.01
86) Perylene-d12	23.797	264	826853	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	953077	116.402	ng	0.00
7) Phenol-d6	6.975	99	1180555	111.233	ng	0.00
23) Nitrobenzene-d5	8.969	82	858739	79.100	ng	0.00
42) 2,4,6-Tribromophenol	15.962	330	779129	155.530	ng	0.00
45) 2-Fluorobiphenyl	13.074	172	2045601	72.963	ng	0.00
79) Terphenyl-d14	19.862	244	4654232	97.451	ng	0.00
Target Compounds						
32) Benzoic acid	9.898	122	36981m	12.703	ng	Qvalue

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

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