

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080423\
 Data File : BM041289.D
 Acq On : 04 Aug 2023 11:51
 Operator : MA/JU
 Sample : 03877-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COMP-1

Manual Integrations
 APPROVED

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

Quant Time: Aug 04 16:13:48 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	100392	20.000	ng	0.00
21) Naphthalene-d8	10.598	136	412351	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	285878	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	765839	20.000	ng	# 0.00
76) Chrysene-d12	21.427	240	752154	20.000	ng	0.00
86) Perylene-d12	23.797	264	715816	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	764871	113.870	ng	0.00
7) Phenol-d6	6.975	99	921160	105.797	ng	0.00
23) Nitrobenzene-d5	8.969	82	704370	79.438	ng	0.00
42) 2,4,6-Tribromophenol	15.962	330	638681	157.949	ng	0.00
45) 2-Fluorobiphenyl	13.074	172	1711044	75.608	ng	0.00
79) Terphenyl-d14	19.862	244	4065824	97.736	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	9.904	122	54210m	17.113	ng	

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080423\
Data File : BM041289.D
Acq On : 04 Aug 2023 11:51
Operator : MA/JU
Sample : 03877-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COMP-1

Quant Time: Aug 04 16:13:48 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

Manual Integrations
APPROVED

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

