

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081325\
 Data File : BF143373.D
 Acq On : 13 Aug 2025 14:11
 Operator : RC/JU
 Sample : Q2827-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-9

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 08/14/2025
 Supervised By :Jagrut Upadhyay 08/15/2025

Quant Time: Aug 13 14:32:25 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 12 13:25:39 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.928	152	174108	20.000	ng	0.00
21) Naphthalene-d8	8.210	136	670714	20.000	ng	0.00
39) Acenaphthene-d10	9.963	164	379214	20.000	ng	0.00
64) Phenanthrene-d10	11.451	188	649011	20.000	ng	0.00
76) Chrysene-d12	14.086	240	441865	20.000	ng	0.00
86) Perylene-d12	15.580	264	377092	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	1181434	117.341	ng	0.02
7) Phenol-d6	6.563	99	1380750	106.940	ng	0.00
23) Nitrobenzene-d5	7.487	82	1056561	98.098	ng	0.00
42) 2,4,6-Tribromophenol	10.757	330	488018	176.092	ng	0.00
45) 2-Fluorobiphenyl	9.281	172	1872645	83.740	ng	0.00
79) Terphenyl-d14	13.033	244	2358587	95.298	ng	0.00
Target Compounds						
32) Benzoic acid	7.969	122	13038m	7.395	ng	Qvalue

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081325\
Data File : BF143373.D
Acq On : 13 Aug 2025 14:11
Operator : RC/JU
Sample : Q2827-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-9

Quant Time: Aug 13 14:32:25 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 12 13:25:39 2025
Response via : Initial Calibration

Manual Integrations
APPROVED

Reviewed By :Rahul Chavli 08/14/2025
Supervised By :Jagrut Upadhyay 08/15/2025

