

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081825\
 Data File : BF143434.D
 Acq On : 18 Aug 2025 11:41
 Operator : RC/JU
 Sample : Q2832-10
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TG-S05

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 08/19/2025
 Supervised By :Jagrut Upadhyay 08/19/2025

Quant Time: Aug 18 12:23:11 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 18 04:40:19 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.928	152	115498	20.000	ng	0.00
21) Naphthalene-d8	8.204	136	444901	20.000	ng	0.00
39) Acenaphthene-d10	9.957	164	246416	20.000	ng	0.00
64) Phenanthrene-d10	11.445	188	401400	20.000	ng	0.00
76) Chrysene-d12	14.086	240	273750	20.000	ng	0.00
86) Perylene-d12	15.574	264	259673	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.569	112	871355	131.145	ng	0.01
7) Phenol-d6	6.557	99	1039114	124.099	ng	0.00
23) Nitrobenzene-d5	7.481	82	765549	102.106	ng	0.00
42) 2,4,6-Tribromophenol	10.751	330	345175	157.378	ng	0.00
45) 2-Fluorobiphenyl	9.280	172	1382683	94.379	ng	0.00
79) Terphenyl-d14	13.027	244	1564193	95.986	ng	0.00
Target Compounds						
32) Benzoic acid	7.975	122	11065m	3.464	ng	Qvalue

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

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