

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093025\
 Data File : BF143844.D
 Acq On : 30 Sep 2025 20:33
 Operator : RC/JU
 Sample : Q3231-08
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-2

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 10/01/2025
 Supervised By :Jagrut Upadhyay 10/01/2025

Quant Time: Oct 01 02:36:40 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF092325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 24 05:00:00 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.886	152	119755	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	445493	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	232539	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	373873	20.000	ng	0.00
76) Chrysene-d12	14.051	240	318319	20.000	ng	-0.01
86) Perylene-d12	15.533	264	494893	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	1061342	143.991	ng	0.01
7) Phenol-d6	6.504	99	1200828	140.999	ng	0.00
23) Nitrobenzene-d5	7.445	82	777602	106.844	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	596664	173.292	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	1575774	105.745	ng	0.00
79) Terphenyl-d14	12.998	244	2027553	129.541	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	7.863	122	13704m	4.481	ng	

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

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