

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080725\
 Data File : BP025315.D
 Acq On : 07 Aug 2025 15:50
 Operator : CG/JU
 Sample : Q2759-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :

BNA_P

ClientSampleId :

SOLID-DRUM

Manual Integrations**APPROVED**

Reviewed By :Rahul Chavli 08/08/2025

Supervised By :Jagrut Upadhyay 08/11/2025

Quant Time: Aug 07 16:08:56 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 06 06:41:44 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	Qvalue
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.807	152	227123	20.000	ng	0.00	
21) Naphthalene-d8	10.578	136	950262	20.000	ng	0.00	
39) Acenaphthene-d10	14.419	164	611259	20.000	ng	0.00	
64) Phenanthrene-d10	17.213	188	1274549	20.000	ng	-0.01	
76) Chrysene-d12	21.660	240	1216767	20.000	ng	0.00	
86) Perylene-d12	25.030	264	1319816	20.000	ng	0.00	
System Monitoring Compounds							
5) 2-Fluorophenol	5.419	112	1515394	107.312	ng	0.01	
7) Phenol-d6	6.972	99	1907321	102.628	ng	0.00	
23) Nitrobenzene-d5	8.949	82	1362404	79.429	ng	0.00	
42) 2,4,6-Tribromophenol	15.931	330	898441	147.333	ng	0.00	
45) 2-Fluorobiphenyl	13.037	172	3119669	70.492	ng	0.00	
79) Terphenyl-d14	19.942	244	3273110	50.808	ng	0.00	
Target Compounds							
10) Phenol	7.002	94	614727	30.914	ng		100
20) 3+4-Methylphenols	8.560	107	5074076	276.939	ng		90
32) Benzoic acid	9.825	122	109401m	18.776	ng		

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

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