

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060925\
 Data File : BP024877.D
 Acq On : 09 Jun 2025 14:52
 Operator : RC/JU
 Sample : Q2230-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GW-MW01-060425

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 06/10/2025
 Supervised By :Jagrut Upadhyay 06/10/2025

Quant Time: Jun 09 15:55:56 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 06 16:20:27 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.608	152	193111	20.000	ng	0.00
21) Naphthalene-d8	10.372	136	768427	20.000	ng	0.00
39) Acenaphthene-d10	14.243	164	479578	20.000	ng	0.00
64) Phenanthrene-d10	17.042	188	943300	20.000	ng	-0.02
76) Chrysene-d12	21.472	240	1244702	20.000	ng	-0.01
86) Perylene-d12	24.713	264	1717483	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.237	112	953456	82.414	ng	0.00
7) Phenol-d6	6.814	99	781439	51.051	ng	0.00
23) Nitrobenzene-d5	8.755	82	1471313	93.041	ng	0.00
42) 2,4,6-Tribromophenol	15.766	330	1067678	161.027	ng	-0.02
45) 2-Fluorobiphenyl	12.854	172	3058686	85.918	ng	0.00
79) Terphenyl-d14	19.772	244	5454769	78.543	ng	-0.02
Target Compounds						
						Qvalue
20) 3+4-Methylphenols	8.408	107	60786	4.043	ng	85
31) Naphthalene	10.425	128	79849	2.028	ng	# 90
32) Benzoic acid	9.713	122	37128	4.631	ng	97
38) 1-Methylnaphthalene	12.266	142	55989m	2.097	ng	

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

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