

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072425\
 Data File : BP025245.D
 Acq On : 24 Jul 2025 16:50
 Operator : CG/JU
 Sample : Q2481-13 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CC0627-CLOXPL

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 07/25/2025
 Supervised By :Jagrut Upadhyay 07/25/2025

Quant Time: Jul 24 17:28:43 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 22 02:55:21 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.437	152	459498	20.000	ng	0.00
21) Naphthalene-d8	10.178	136	1166103	20.000	ng	# 0.00
39) Acenaphthene-d10	14.095	164	4344m	20.000	ng	0.02
64) Phenanthrene-d10	16.889	188	1393965	20.000	ng	# 0.00
76) Chrysene-d12	21.324	240	1310534	20.000	ng	# 0.00
86) Perylene-d12	24.471	264	8832	20.000	ng	# 0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.131	112	60381	2.220	ng	0.03
7) Phenol-d6	0.000	99	0d	0.000	ng	
23) Nitrobenzene-d5	8.566	82	329335	15.695	ng	0.00
42) 2,4,6-Tribromophenol	15.607	330	197724	3087.501	ng	0.00
45) 2-Fluorobiphenyl	12.684	172	786313	2551.923	ng	0.00
79) Terphenyl-d14	19.654	244	1050314	15.821	ng	0.00
Target Compounds						
25) Isophorone	9.196	82	200737	5.440	ng	# 98

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

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