

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060925\
 Data File : BP024875.D
 Acq On : 09 Jun 2025 13:31
 Operator : RC/JU
 Sample : Q2209-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P01W

Quant Time: Jun 09 13:51:31 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	305772	20.000	ng	0.00
21) Naphthalene-d8	10.378	136	1202322	20.000	ng	0.00
39) Acenaphthene-d10	14.248	164	730596	20.000	ng	0.00
64) Phenanthrene-d10	17.048	188	1469775	20.000	ng	-0.01
76) Chrysene-d12	21.466	240	1832304	20.000	ng	-0.02
86) Perylene-d12	24.712	264	2370307	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.237	112	1664555	90.868	ng	0.00
7) Phenol-d6	6.813	99	1441905	59.491	ng	0.00
23) Nitrobenzene-d5	8.760	82	2205524	89.138	ng	0.00
42) 2,4,6-Tribromophenol	15.766	330	1613344	159.723	ng	-0.02
45) 2-Fluorobiphenyl	12.854	172	4535208	83.623	ng	0.00
79) Terphenyl-d14	19.772	244	8128028	79.503	ng	-0.02

Target Compounds	Qvalue

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

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