

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061325\
 Data File : BP024939.D
 Acq On : 13 Jun 2025 11:23
 Operator : RC/JU
 Sample : PB168437BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB168437BL

Quant Time: Jun 13 12:07:39 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	274334	20.000	ng	0.00
21) Naphthalene-d8	10.378	136	1046404	20.000	ng	0.00
39) Acenaphthene-d10	14.254	164	653672	20.000	ng	0.00
64) Phenanthrene-d10	17.054	188	1304490	20.000	ng	0.00
76) Chrysene-d12	21.489	240	1274406	20.000	ng	0.00
86) Perylene-d12	24.754	264	1474268	20.000	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.243	112	2551476	155.246	ng	0.00
7) Phenol-d6	6.819	99	3124700	143.696	ng	0.00
23) Nitrobenzene-d5	8.760	82	1934875	89.852	ng	0.00
42) 2,4,6-Tribromophenol	15.778	330	1546217	171.091	ng	0.00
45) 2-Fluorobiphenyl	12.860	172	4514889	93.045	ng	0.00
79) Terphenyl-d14	19.789	244	6901924	97.065	ng	0.00

Target Compounds	Qvalue

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

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