

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022825\
 Data File : BP024133.D
 Acq On : 28 Feb 2025 13:31
 Operator : RC/JU
 Sample : PB166890BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB166890BL

Quant Time: Feb 28 13:52:41 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP021725.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 18 02:25:07 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	367881	20.000	ng	0.00
21) Naphthalene-d8	10.540	136	1405133	20.000	ng	0.00
39) Acenaphthene-d10	14.393	164	835299	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	1671238	20.000	ng	0.00
76) Chrysene-d12	21.651	240	1717685	20.000	ng	0.00
86) Perylene-d12	25.051	264	1892079	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	3469344	147.152	ng	0.00
7) Phenol-d6	6.946	99	4265874	140.777	ng	0.00
23) Nitrobenzene-d5	8.910	82	2488278	98.794	ng	0.00
42) 2,4,6-Tribromophenol	15.910	330	1708840	167.018	ng	0.00
45) 2-Fluorobiphenyl	12.998	172	5550897	92.903	ng	0.00
79) Terphenyl-d14	19.922	244	9050146	104.256	ng	0.00

Target Compounds	Qvalue

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

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