

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021725\
 Data File : BP024078.D
 Acq On : 17 Feb 2025 17:36
 Operator : RC/JU
 Sample : PB166698BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB166698BL

Quant Time: Feb 18 03:02:42 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	7.764	152	437383	20.000	ng	0.00
21) Naphthalene-d8	10.534	136	1680499	20.000	ng	0.00
39) Acenaphthene-d10	14.393	164	989916	20.000	ng	0.00
64) Phenanthrene-d10	17.187	188	1884903	20.000	ng	0.00
76) Chrysene-d12	21.633	240	1509617	20.000	ng	-0.01
86) Perylene-d12	25.016	264	1468519	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	4325761	154.321	ng	0.00
7) Phenol-d6	6.946	99	5278716	146.520	ng	0.00
23) Nitrobenzene-d5	8.911	82	3043370	101.033	ng	0.00
42) 2,4,6-Tribromophenol	15.904	330	1933104	159.426	ng	0.00
45) 2-Fluorobiphenyl	13.004	172	7140735	100.844	ng	0.00
79) Terphenyl-d14	19.910	244	9686972	126.973	ng	0.00

Target Compounds	Qvalue

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

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