

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060324\
 Data File : BP020488.D
 Acq On : 03 Jun 2024 11:56
 Operator : MA/JU
 Sample : PB161236BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB161236BL

Quant Time: Jun 03 13:09:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 30 03:41:03 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	334266	20.000	ng	0.00
21) Naphthalene-d8	10.528	136	1295439	20.000	ng	0.00
39) Acenaphthene-d10	14.387	164	826371	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	1615176	20.000	ng	-0.01
76) Chrysene-d12	21.651	240	1107913	20.000	ng	-0.02
86) Perylene-d12	25.004	264	1162256	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	2648360	141.856	ng	0.00
7) Phenol-d6	6.928	99	3385914	128.603	ng	0.00
23) Nitrobenzene-d5	8.910	82	2121317	89.638	ng	-0.01
42) 2,4,6-Tribromophenol	15.904	330	1106566	129.029	ng	0.00
45) 2-Fluorobiphenyl	12.998	172	4914319	88.652	ng	0.00
79) Terphenyl-d14	19.933	244	6160949	92.573	ng	-0.01

Target Compounds	Qvalue

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

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