

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092324\
Data File : BP021977.D
Acq On : 23 Sep 2024 11:09
Operator : RC/JU
Sample : PB163139BL
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB163139BL

Quant Time: Sep 23 11:40:06 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 20 01:22:00 2024
Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	74719	20.000	ng	0.00
21) Naphthalene-d8	10.469	136	253893	20.000	ng	0.00
39) Acenaphthene-d10	14.322	164	170443	20.000	ng	0.00
64) Phenanthrene-d10	17.116	188	380108	20.000	ng	-0.02
76) Chrysene-d12	21.551	240	434852	20.000	ng	#-0.01
86) Perylene-d12	24.839	264	535492	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.334	112	487019	122.491	ng	0.00
7) Phenol-d6	6.893	99	589392	114.190	ng	0.00
23) Nitrobenzene-d5	8.852	82	399997	83.363	ng	0.00
42) 2,4,6-Tribromophenol	15.834	330	339252	107.364	ng	-0.02
45) 2-Fluorobiphenyl	12.934	172	966303	82.759	ng	0.00
79) Terphenyl-d14	19.851	244	1980260	85.544	ng	0.00

Target Compounds	Qvalue

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

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