

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092024\
Data File : BP021972.D
Acq On : 20 Sep 2024 18:13
Operator : RC/JU
Sample : PB163110BL
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB163110BL

Quant Time: Sep 21 01:58:40 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	104663	20.000	ng	0.00
21) Naphthalene-d8	10.475	136	353857	20.000	ng	0.00
39) Acenaphthene-d10	14.333	164	228261	20.000	ng	0.00
64) Phenanthrene-d10	17.122	188	496320	20.000	ng	-0.02
76) Chrysene-d12	21.557	240	510780	20.000	ng	0.00
86) Perylene-d12	24.851	264	583237	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.334	112	753100	135.223	ng	0.00
7) Phenol-d6	6.893	99	936678	129.554	ng	0.00
23) Nitrobenzene-d5	8.851	82	635147	94.975	ng	0.00
42) 2,4,6-Tribromophenol	15.833	330	529238	125.065	ng	-0.02
45) 2-Fluorobiphenyl	12.939	172	1525661	97.569	ng	0.00
79) Terphenyl-d14	19.851	244	2812027	103.417	ng	0.00

Target Compounds	Qvalue

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

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