

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092324\
 Data File : BP021982.D
 Acq On : 23 Sep 2024 14:32
 Operator : RC/JU
 Sample : PB163139BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB163139BL

Quant Time: Sep 23 14:57:54 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	7.711	152	98936	20.000	ng	0.00
21) Naphthalene-d8	10.475	136	342173	20.000	ng	0.00
39) Acenaphthene-d10	14.328	164	225444	20.000	ng	0.00
64) Phenanthrene-d10	17.122	188	541145	20.000	ng	-0.02
76) Chrysene-d12	21.539	240	653961	20.000	ng	-0.02
86) Perylene-d12	24.827	264	768319	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.334	112	737122	140.015	ng	0.00
7) Phenol-d6	6.893	99	896699	131.204	ng	0.00
23) Nitrobenzene-d5	8.852	82	612402	94.701	ng	0.00
42) 2,4,6-Tribromophenol	15.834	330	570144	136.415	ng	-0.02
45) 2-Fluorobiphenyl	12.940	172	1481715	95.942	ng	0.00
79) Terphenyl-d14	19.845	244	3403802	97.773	ng	-0.01

Target Compounds	Qvalue

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

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