

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110824\
Data File : BP022937.D
Acq On : 08 Nov 2024 16:51
Operator : RC/JU
Sample : PB164369BL
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB164369BL

Quant Time: Nov 08 17:47:15 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110724.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 07 23:36:11 2024
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	60434	20.000	ng	0.00
21) Naphthalene-d8	10.351	136	205441	20.000	ng	0.00
39) Acenaphthene-d10	14.216	164	151785	20.000	ng	0.00
64) Phenanthrene-d10	17.033	188	366674	20.000	ng	0.01
76) Chrysene-d12	21.468	240	471057	20.000	ng	0.00
86) Perylene-d12	24.704	264	585630	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.234	112	465033	146.032	ng	0.00
7) Phenol-d6	6.793	99	589166	132.666	ng	0.00
23) Nitrobenzene-d5	8.740	82	414084	95.224	ng	0.00
42) 2,4,6-Tribromophenol	15.728	330	396493	134.767	ng	-0.02
45) 2-Fluorobiphenyl	12.816	172	1030942	97.050	ng	0.00
79) Terphenyl-d14	19.757	244	2519361	100.442	ng	0.00

Target Compounds	Qvalue

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

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