

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042325\
 Data File : BP024394.D
 Acq On : 23 Apr 2025 13:27
 Operator : RC/JU
 Sample : PB167660TB
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB167660TB

Quant Time: Apr 23 13:56:58 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 18 12:04:48 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	223546	20.000	ng	0.00
21) Naphthalene-d8	10.487	136	891166	20.000	ng	-0.01
39) Acenaphthene-d10	14.345	164	538193	20.000	ng	0.00
64) Phenanthrene-d10	17.151	188	1025007	20.000	ng	0.00
76) Chrysene-d12	21.592	240	1045807	20.000	ng	0.00
86) Perylene-d12	24.939	264	1193660	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.346	112	1920255	142.031	ng	0.00
7) Phenol-d6	6.905	99	2362290	127.605	ng	0.00
23) Nitrobenzene-d5	8.869	82	1431310	91.582	ng	0.00
42) 2,4,6-Tribromophenol	15.857	330	1085213	145.741	ng	0.00
45) 2-Fluorobiphenyl	12.957	172	3262478	92.815	ng	0.00
79) Terphenyl-d14	19.875	244	5014982	96.656	ng	0.00

Target Compounds	Qvalue

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

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