

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110724\
 Data File : BP022923.D
 Acq On : 07 Nov 2024 18:15
 Operator : RC/JU
 Sample : PB164711BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB164711BL

Quant Time: Nov 08 00:03:14 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	59975	20.000	ng	0.00
21) Naphthalene-d8	10.346	136	207655	20.000	ng	0.00
39) Acenaphthene-d10	14.222	164	160890	20.000	ng	0.00
64) Phenanthrene-d10	17.028	188	390642	20.000	ng	0.00
76) Chrysene-d12	21.457	240	463436	20.000	ng	0.00
86) Perylene-d12	24.668	264	541811	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.234	112	539258	170.636	ng	0.00
7) Phenol-d6	6.793	99	688378	156.192	ng	0.00
23) Nitrobenzene-d5	8.734	82	481123	109.461	ng	0.00
42) 2,4,6-Tribromophenol	15.745	330	538427	172.653	ng	0.00
45) 2-Fluorobiphenyl	12.822	172	1247140	110.758	ng	0.00
79) Terphenyl-d14	19.757	244	2911137	117.970	ng	0.00

Target Compounds	Qvalue

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

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