

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112322\
 Data File : BP012705.D
 Acq On : 23 Nov 2022 11:06
 Operator : CG/JU
 Sample : PB149242BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB149242BL

Quant Time: Nov 24 02:25:56 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 00:38:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	325711	20.000	ng	0.00
21) Naphthalene-d8	10.851	136	1377434	20.000	ng	0.00
39) Acenaphthene-d10	14.669	164	861573	20.000	ng	0.00
64) Phenanthrene-d10	17.428	188	1845452	20.000	ng	0.00
76) Chrysene-d12	21.510	240	1582993	20.000	ng	0.00
86) Perylene-d12	24.039	264	1341732	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	2744605	182.555	ng	0.00
7) Phenol-d6	7.187	99	3637097	168.687	ng	0.00
23) Nitrobenzene-d5	9.199	82	2336207	94.590	ng	0.00
42) 2,4,6-Tribromophenol	16.157	330	1331479	120.633	ng	0.00
45) 2-Fluorobiphenyl	13.298	172	5800575	95.916	ng	0.00
79) Terphenyl-d14	19.974	244	9272263	124.492	ng	0.00

Target Compounds	Qvalue

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

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