

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112222\
 Data File : BP012702.D
 Acq On : 22 Nov 2022 17:20
 Operator : CG/JU
 Sample : PB149238BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB149238BL

Quant Time: Nov 23 06:25:07 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	8.040	152	298111	20.000	ng	0.00
21) Naphthalene-d8	10.858	136	1213022	20.000	ng	0.00
39) Acenaphthene-d10	14.675	164	777258	20.000	ng	0.00
64) Phenanthrene-d10	17.434	188	1704387	20.000	ng	0.00
76) Chrysene-d12	21.522	240	1555997	20.000	ng	0.01
86) Perylene-d12	24.051	264	1409360	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	2514580	182.740	ng	0.00
7) Phenol-d6	7.187	99	3358600	170.192	ng	0.00
23) Nitrobenzene-d5	9.205	82	2123588	97.635	ng	0.00
42) 2,4,6-Tribromophenol	16.169	330	1351530	134.130	ng	0.01
45) 2-Fluorobiphenyl	13.304	172	5389665	98.789	ng	0.00
79) Terphenyl-d14	19.986	244	9026876	123.300	ng	0.01

Target Compounds Qvalue

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

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