

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111222\
 Data File : BP012632.D
 Acq On : 12 Nov 2022 22:31
 Operator : CG/JU
 Sample : N5502-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 2182RE

Quant Time: Nov 14 03:23:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Nov 13 22:44:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.711	152	299881	20.000	ng	0.00
21) Naphthalene-d8	10.510	136	1221918	20.000	ng	0.00
39) Acenaphthene-d10	14.369	164	801050	20.000	ng	-0.01
64) Phenanthrene-d10	17.134	188	1747896	20.000	ng	-0.01
76) Chrysene-d12	21.233	240	1596414	20.000	ng	-0.01
86) Perylene-d12	23.574	264	1373937	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.299	112	2503328	155.440	ng	0.00
7) Phenol-d6	6.905	99	3033214	135.689	ng	0.00
23) Nitrobenzene-d5	8.887	82	2183698	99.146	ng	0.00
42) 2,4,6-Tribromophenol	15.875	330	1819579	166.393	ng	-0.01
45) 2-Fluorobiphenyl	12.987	172	4972459	94.234	ng	-0.01
79) Terphenyl-d14	19.710	244	7741897	98.248	ng	0.00

Target Compounds Qvalue

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

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