

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010322\
 Data File : BP008641.D
 Acq On : 03 Jan 2022 17:07
 Operator : CG/JU
 Sample : M5257-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P3-B

Quant Time: Jan 04 05:59:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP122821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 30 08:44:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	552338	20.000	ng	-0.01
21) Naphthalene-d8	10.693	136	1987234	20.000	ng	0.00
39) Acenaphthene-d10	14.522	164	1180675	20.000	ng	0.00
64) Phenanthrene-d10	17.269	188	2309313	20.000	ng	0.00
76) Chrysene-d12	21.351	240	2422954	20.000	ng	-0.01
86) Perylene-d12	23.774	264	2811578	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.470	112	3534835	116.813	ng	0.00
7) Phenol-d6	7.058	99	3842865	91.236	ng	0.00
23) Nitrobenzene-d5	9.046	82	2813351	84.219	ng	-0.01
42) 2,4,6-Tribromophenol	16.010	330	2252020	140.980	ng	0.00
45) 2-Fluorobiphenyl	13.146	172	7410705	90.938	ng	-0.01
79) Terphenyl-d14	19.828	244	10541138	88.829	ng	0.00

Target Compounds	Qvalue

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

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