

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010322\
 Data File : BP008648.D
 Acq On : 03 Jan 2022 21:04
 Operator : CG/JU
 Sample : M5257-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P2-B

Quant Time: Jan 04 06:01:44 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	576222	20.000	ng	-0.01
21) Naphthalene-d8	10.687	136	2117202	20.000	ng	-0.01
39) Acenaphthene-d10	14.516	164	1244179	20.000	ng	-0.01
64) Phenanthrene-d10	17.269	188	2341554	20.000	ng	0.00
76) Chrysene-d12	21.351	240	2249863	20.000	ng	-0.01
86) Perylene-d12	23.768	264	2470356	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	4160297	131.784	ng	0.00
7) Phenol-d6	7.058	99	4555078	103.663	ng	0.00
23) Nitrobenzene-d5	9.046	82	3377966	94.913	ng	-0.01
42) 2,4,6-Tribromophenol	16.010	330	2262624	134.414	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	8389221	97.691	ng	-0.01
79) Terphenyl-d14	19.827	244	10742659	97.491	ng	0.00
Target Compounds						
32) Benzoic acid	9.934	122	69311	3.777	ng	Qvalue 97

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

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