

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010322\
 Data File : BP008647.D
 Acq On : 03 Jan 2022 20:30
 Operator : CG/JU
 Sample : M5257-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 P2-A

Quant Time: Jan 04 06:01:32 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	609074	20.000	ng	-0.01
21) Naphthalene-d8	10.687	136	2309070	20.000	ng	-0.01
39) Acenaphthene-d10	14.516	164	1434152	20.000	ng	-0.01
64) Phenanthrene-d10	17.269	188	2813196	20.000	ng	0.00
76) Chrysene-d12	21.351	240	2516438	20.000	ng	-0.01
86) Perylene-d12	23.774	264	2687384	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	4467468	133.881	ng	0.00
7) Phenol-d6	7.058	99	4884464	105.163	ng	0.00
23) Nitrobenzene-d5	9.046	82	3820944	98.439	ng	-0.01
42) 2,4,6-Tribromophenol	16.010	330	2853253	147.048	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	9558308	96.561	ng	-0.01
79) Terphenyl-d14	19.827	244	12169215	98.739	ng	0.00
Target Compounds						
32) Benzoic acid	9.940	122	87963	4.396	ng	94
50) Dimethylphthalate	13.975	163	224758	2.035	ng	100

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

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