

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102922\
 Data File : BM037272.D
 Acq On : 29 Oct 2022 15:52
 Operator : CG/JU
 Sample : N5236-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LSP-GW-02

Quant Time: Oct 31 02:06:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 31 01:34:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	469894	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	1847818	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	1010881	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1729959	20.000	ng	0.00
76) Chrysene-d12	21.421	240	1154159	20.000	ng	0.00
86) Perylene-d12	23.780	264	1295376	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	3125291	114.906	ng	0.00
7) Phenol-d6	7.046	99	2915361	78.976	ng	0.00
23) Nitrobenzene-d5	9.045	82	4448264	121.455	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	2328442	195.456	ng	0.00
45) 2-Fluorobiphenyl	13.133	172	9409390	123.676	ng	0.00
79) Terphenyl-d14	19.868	244	9768138	151.772	ng	0.00

Target Compounds	Qvalue

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

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