

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040522\
 Data File : BP009694.D
 Acq On : 05 Apr 2022 12:47
 Operator : CG/JU
 Sample : N2249-01DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-1RDL

Quant Time: Apr 05 13:41:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 05 13:35:48 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	238549	20.000	ng	0.00
21) Naphthalene-d8	10.516	136	940313	20.000	ng	0.00
39) Acenaphthene-d10	14.375	164	601630	20.000	ng	0.00
64) Phenanthrene-d10	17.139	188	1289186	20.000	ng	0.00
76) Chrysene-d12	21.263	240	1372251	20.000	ng	0.00
86) Perylene-d12	23.633	264	1446423	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	172183	12.602	ng	0.01
7) Phenol-d6	6.928	99	186394	10.088	ng	0.01
23) Nitrobenzene-d5	8.887	82	239330	13.525	ng	0.00
42) 2,4,6-Tribromophenol	15.881	330	148056	14.913	ng	0.00
45) 2-Fluorobiphenyl	12.993	172	601507	13.860	ng	0.00
79) Terphenyl-d14	19.722	244	761593	9.962	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.563	128	2170887	44.820	ng	100
37) 2-Methylnaphthalene	12.187	142	262814	7.733	ng	100
38) 1-Methylnaphthalene	12.404	142	153723	4.643	ng	99

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

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