

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120621\
 Data File : BP008237.D
 Acq On : 06 Dec 2021 19:52
 Operator : CG/JU
 Sample : M4917-07DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-23DL

Quant Time: Dec 07 05:11:47 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 03 14:39:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	65876	20.000	ng	-0.01
21) Naphthalene-d8	10.587	136	247931	20.000	ng	-0.01
39) Acenaphthene-d10	14.434	164	164269	20.000	ng	-0.01
64) Phenanthrene-d10	17.198	188	375727	20.000	ng	-0.01
76) Chrysene-d12	21.304	240	436307	20.000	ng	-0.01
86) Perylene-d12	23.698	264	489226	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	45614	11.149	ng	0.00
7) Phenol-d6	6.981	99	34422	6.232	ng	0.00
23) Nitrobenzene-d5	8.946	82	80088	14.689	ng	-0.02
42) 2,4,6-Tribromophenol	15.940	330	64619	30.776	ng	-0.01
45) 2-Fluorobiphenyl	13.057	172	190598	16.853	ng	-0.01
79) Terphenyl-d14	19.769	244	332260	15.915	ng	-0.01
Target Compounds						
						Qvalue
31) Naphthalene	10.634	128	705336	55.535	ng	99
37) 2-Methylnaphthalene	12.251	142	173886	19.358	ng	98
38) 1-Methylnaphthalene	12.469	142	130788	14.958	ng	98

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

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