

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112420\
 Data File : BP004084.D
 Acq On : 24 Nov 2020 19:51
 Operator : CG/JU
 Sample : L4829-07DL2 25X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-12DL2

Quant Time: Nov 24 23:59:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 23 16:42:41 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.269	152	112478	20.00	ng	0.00
21) Naphthalene-d8	11.099	136	436307	20.00	ng	# 0.00
39) Acenaphthene-d10	14.887	164	274663	20.00	ng	0.00
64) Phenanthrene-d10	17.622	188	586612	20.00	ng	# 0.00
76) Chrysene-d12	21.657	240	563724	20.00	ng	0.00
86) Perylene-d12	24.127	264	601961	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.805	112	12468	1.85	ng	0.00
7) Phenol-d6	7.458	99	9308	0.96	ng	0.00
23) Nitrobenzene-d5	9.434	82	25815	2.90	ng	0.00
42) 2,4,6-Tribromophenol	16.375	330	15360	4.47	ng	0.00
45) 2-Fluorobiphenyl	13.510	172	63760	3.25	ng	0.00
79) Terphenyl-d14	20.116	244	78586	2.69	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	11.146	128	1562857	66.75	ng	100
37) 2-Methylnaphthalene	12.734	142	193396	11.54	ng	99
38) 1-Methylnaphthalene	12.951	142	339556	21.25	ng	100
52) Acenaphthene	14.951	154	63797	3.73	ng	99
71) Phenanthrene	17.663	178	101084	3.07	ng	98

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

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