

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121721\
 Data File : BP008486.D
 Acq On : 17 Dec 2021 23:32
 Operator : CG/JU
 Sample : M4873-07DL 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A2-8-6.5DL

Quant Time: Dec 18 07:48:56 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP121421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 18 06:49:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.728	152	61535	20.000	ng	0.00
21) Naphthalene-d8	10.522	136	228592	20.000	ng	0.00
39) Acenaphthene-d10	14.380	164	150691	20.000	ng	0.00
64) Phenanthrene-d10	17.145	188	320921	20.000	ng	0.00
76) Chrysene-d12	21.257	240	368222	20.000	ng	0.00
86) Perylene-d12	23.627	264	404368	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.346	112	26393	6.774	ng	0.02
7) Phenol-d6	6.946	99	36487	6.972	ng	0.02
23) Nitrobenzene-d5	8.898	82	28499	5.282	ng	0.00
42) 2,4,6-Tribromophenol	15.904	330	11148	5.034	ng	0.01
45) 2-Fluorobiphenyl	12.998	172	63447	5.352	ng	0.00
79) Terphenyl-d14	19.721	244	107890	5.099	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.569	128	453393	37.822	ng	97
37) 2-Methylnaphthalene	12.192	142	442378	53.543	ng	98
38) 1-Methylnaphthalene	12.410	142	214632	26.683	ng	99
46) 1,1'-Biphenyl	13.210	154	32766	2.814	ng	99
71) Phenanthrene	17.192	178	67397	3.877	ng	97

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

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