

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120721\
 Data File : BP008271.D
 Acq On : 08 Dec 2021 02:51
 Operator : CG/JU
 Sample : M4932-01DL 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DRUM1-2DL

Quant Time: Dec 08 03:50:37 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	7.793	152	67928	20.000	ng	0.00
21) Naphthalene-d8	10.587	136	266366	20.000	ng	#-0.01
39) Acenaphthene-d10	14.434	164	192208	20.000	ng	-0.01
64) Phenanthrene-d10	17.198	188	427104	20.000	ng	-0.01
76) Chrysene-d12	21.304	240	468976	20.000	ng	-0.01
86) Perylene-d12	23.698	264	513176	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	64616	15.316	ng	0.00
7) Phenol-d6	6.981	99	94083	16.520	ng	0.00
23) Nitrobenzene-d5	8.951	82	129215	22.059	ng	-0.01
42) 2,4,6-Tribromophenol	15.933	330	39841	16.217	ng	-0.02
45) 2-Fluorobiphenyl	13.051	172	113531	8.580	ng	-0.02
79) Terphenyl-d14	19.769	244	195601	8.717	ng	-0.01
Target Compounds						
						Qvalue
27) 2,4-Dimethylphenol	9.804	122	30855	8.418	ng	# 66
31) Naphthalene	10.640	128	533022	39.063	ng	# 94
37) 2-Methylnaphthalene	12.251	142	22952	2.378	ng	# 92

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

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