

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120321\
 Data File : BP008208.D
 Acq On : 03 Dec 2021 16:53
 Operator : CG/JU
 Sample : M4798-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-3-4.5

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/06/2021
 Supervised By :Sohil Jodhani 12/06/2021

Quant Time: Dec 03 17:27:19 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	76744	20.000	ng	0.00
21) Naphthalene-d8	10.593	136	287005	20.000	ng	# 0.00
39) Acenaphthene-d10	14.451	164	187822	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	412507	20.000	ng	0.00
76) Chrysene-d12	21.310	240	500791	20.000	ng	0.00
86) Perylene-d12	23.704	264	484211	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	558611	117.197	ng	0.00
7) Phenol-d6	6.981	99	718325	111.639	ng	0.00
23) Nitrobenzene-d5	8.957	82	500011	79.222	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	315336	131.353	ng	0.00
45) 2-Fluorobiphenyl	13.069	172	976566	75.523	ng	0.00
79) Terphenyl-d14	19.780	244	1873691	78.194	ng	0.00
Target Compounds						Qvalue
38) 1-Methylnaphthalene	12.481	142	64046	6.328	ng	# 98
52) Acenaphthene	14.516	154	35127	3.550	ng	# 88
58) Fluorene	15.504	166	47149m	3.432	ng	

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

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