

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120321\
 Data File : BM033287.D
 Acq On : 03 Dec 2021 15:37
 Operator : CG/JU
 Sample : M4798-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-1-4.5

Quant Time: Dec 03 16:08:13 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 24 15:16:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	120753	20.000	ng	0.00
21) Naphthalene-d8	10.725	136	489112	20.000	ng	#-0.01
39) Acenaphthene-d10	14.554	164	311052	20.000	ng	0.00
64) Phenanthrene-d10	17.295	188	629249	20.000	ng	# 0.00
76) Chrysene-d12	21.447	240	618923	20.000	ng	# 0.00
86) Perylene-d12	23.783	264	646849	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.519	112	909522	124.383	ng	0.01
7) Phenol-d6	7.119	99	1206826	123.231	ng	0.02
23) Nitrobenzene-d5	9.089	82	862613	81.874	ng	-0.01
42) 2,4,6-Tribromophenol	16.042	330	358458	105.716	ng	0.00
45) 2-Fluorobiphenyl	13.177	172	1599794	72.534	ng	0.00
79) Terphenyl-d14	19.901	244	2378068	75.977	ng	0.00
Target Compounds						Qvalue
38) 1-Methylnaphthalene	12.601	142	209348	12.216	ng	100
52) Acenaphthene	14.618	154	74992	4.373	ng	95
55) Dibenzofuran	14.954	168	74412	2.609	ng	# 1
58) Fluorene	15.601	166	132226	5.922	ng	# 81
60) Diethylphthalate	15.360	149	66630	3.038	ng	# 66
71) Phenanthrene	17.336	178	207413	5.997	ng	# 81
75) Fluoranthene	19.342	202	99897	2.513	ng	83
78) Pyrene	19.701	202	172056	4.223	ng	# 94

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

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