

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121021\
 Data File : BM033405.D
 Acq On : 10 Dec 2021 21:38
 Operator : CG/JU
 Sample : M4873-08
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A2-9-6.5

Quant Time: Dec 11 02:15:15 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM120821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 09 07:11:07 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.913	152	87377	20.000	ng	0.00
21) Naphthalene-d8	10.713	136	348721	20.000	ng	# 0.00
39) Acenaphthene-d10	14.542	164	218275	20.000	ng	0.00
64) Phenanthrene-d10	17.277	188	363140	20.000	ng	# 0.00
76) Chrysene-d12	21.430	240	314488	20.000	ng	-0.01
86) Perylene-d12	23.747	264	300395	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.507	112	296481	55.665	ng	0.01
7) Phenol-d6	7.090	99	620759	83.588	ng	0.00
23) Nitrobenzene-d5	9.090	82	593972	72.322	ng	0.01
42) 2,4,6-Tribromophenol	16.024	330	129088	55.285	ng	0.00
45) 2-Fluorobiphenyl	13.160	172	905773	57.400	ng	0.00
79) Terphenyl-d14	19.877	244	1030285	60.695	ng	-0.01
Target Compounds						
						Qvalue
10) Phenol	7.113	94	22008	2.969	ng	# 41
31) Naphthalene	10.778	128	8568945	459.782	ng	98
37) 2-Methylnaphthalene	12.383	142	7737758	607.364	ng	98
38) 1-Methylnaphthalene	12.595	142	3813883	303.586	ng	98
46) 1,1'-Biphenyl	13.372	154	517874	31.150	ng	97
52) Acenaphthene	14.601	154	138419	11.381	ng	# 88
55) Dibenzofuran	14.936	168	230124	11.383	ng	# 1
58) Fluorene	15.583	166	355548	22.347	ng	# 88
71) Phenanthrene	17.318	178	967412	47.584	ng	# 96
72) Anthracene	17.413	178	43780	2.204	ng	# 59
78) Pyrene	19.683	202	100504	4.563	ng	# 88

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

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