

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071423\
 Data File : BM040891.D
 Acq On : 14 Jul 2023 22:41
 Operator : MA/JU
 Sample : 03437-10
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A46K7

Quant Time: Jul 15 04:00:26 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 15 03:58:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.861	152	4882	0.400	ng/ul	0.00
4) Naphthalene-d8	10.663	136	16099	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.509	164	6747	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.257	188	12084	0.400	ng/ul	0.00
17) Chrysene-d12	21.448	240	9892	0.400	ng/ul	# 0.00
23) Perylene-d12	23.824	264	9882	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.279	96	26117	3.406	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.258	152	4974	0.220	ng/ul	0.00
18) Fluoranthene-d10	19.292	212	9636	0.304	ng/ul	0.00
Target Compounds						
						Qvalue
5) Naphthalene	10.713	128	595743	13.416	ng/ul	99
8) 1-Methylnaphthalene	12.544	142	482551	18.515	ng/ul	100
11) Acenaphthene	14.569	153	448215	17.609	ng/ul	98
12) Fluorene	15.559	166	6452	0.237	ng/ul#	97
14) Pentachlorophenol	16.911	266	169908	73.024	ng/ul	93
15) Phenanthrene	17.299	178	662102	17.639	ng/ul	99
16) Anthracene	17.392	178	828570	26.268	ng/ul	99
19) Fluoranthene	19.320	202	1972219	46.936	ng/ul	95
20) Pyrene	19.687	202	1353876	30.054	ng/ul	94
21) Benzo(a)anthracene	21.433	228	1557631	46.552	ng/ul	99
22) Chrysene	21.483	228	955033	25.022	ng/ul	99
24) Benzo(b)fluoranthene	23.099	252	1012354	26.471	ng/ul	85
26) Benzo(a)pyrene	23.722	252	884441	25.674	ng/ul#	80
27) Indeno(1,2,3-cd)pyrene	26.290	276	1266310	26.395	ng/ul#	91
28) Dibenzo(a,h)anthracene	26.300	278	802587	21.366	ng/ul#	91

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

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