

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071023\
 Data File : BM040793.D
 Acq On : 10 Jul 2023 10:40
 Operator : MA/JU
 Sample : 03384-03MSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YBW82MSD

Quant Time: Jul 11 03:55:37 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 : Tue Jul 11 03:54:24 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	2634	0.400	ng/ul	0.00
4) Naphthalene-d8	10.691	136	10072	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.532	164	4669	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.286	188	7257	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.477	240	5081	0.400	ng/ul	0.00
23) Perylene-d12	23.868	264	4271	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	1022	0.247	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.280	152	3821	0.271	ng/ul	0.00
18) Fluoranthene-d10	19.315	212	6490	0.398	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	2551	0.606	ng/ul	94
5) Naphthalene	10.741	128	8512	0.306	ng/ul	97
7) 2-Methylnaphthalene	12.357	142	4555	0.279	ng/ul	95
8) 1-Methylnaphthalene	12.572	142	4469	0.274	ng/ul	97
10) Acenaphthylene	14.254	152	6198	0.297	ng/ul	95
11) Acenaphthene	14.592	153	5727	0.325	ng/ul	96
12) Fluorene	15.587	166	5879	0.313	ng/ul	98
14) Pentachlorophenol	16.944	266	530	0.379	ng/ul	91
15) Phenanthrene	17.329	178	7898	0.350	ng/ul	96
16) Anthracene	17.426	178	6996	0.369	ng/ul#	93
19) Fluoranthene	19.343	202	8299	0.385	ng/ul	97
20) Pyrene	19.710	202	8886	0.384	ng/ul	98
21) Benzo(a)anthracene	21.463	228	5754	0.335	ng/ul	98
22) Chrysene	21.515	228	6946	0.354	ng/ul	98
24) Benzo(b)fluoranthene	23.140	252	6203	0.375	ng/ul	89
25) Benzo(k)fluoranthene	23.187	252	6674	0.410	ng/ul#	86
26) Benzo(a)pyrene	23.766	252	5360	0.360	ng/ul#	77
27) Indeno(1,2,3-cd)pyrene	26.354	276	7432	0.358	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.371	278	5849	0.360	ng/ul#	91
29) Benzo(g,h,i)perylene	27.115	276	6566	0.361	ng/ul	94

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

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