

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092223\
 Data File : BM042020.D
 Acq On : 23 Sep 2023 11:49
 Operator : MA/JU
 Sample : PB155506BS
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS506

Quant Time: Sep 25 01:04:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 13:50:58 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.950	152	5479	0.400	ng/ul	0.00
4) Naphthalene-d8	10.763	136	12852	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.597	164	5692	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.350	188	10235	0.400	ng/ul	0.00
17) Chrysene-d12	21.521	240	3815	0.400	ng/ul	0.00
23) Perylene-d12	23.930	264	3838	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	6389	0.856	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.346	152	7538	0.468	ng/ul	0.00
18) Fluoranthene-d10	19.371	212	8600	0.547	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.406	88	15255	1.942	ng/ul#	87
5) Naphthalene	10.818	128	17893	0.416	ng/ul	100
7) 2-Methylnaphthalene	12.418	142	10544	0.411	ng/ul	99
8) 1-Methylnaphthalene	12.643	142	11341	0.435	ng/ul	99
10) Acenaphthylene	14.324	152	15465	0.405	ng/ul	99
11) Acenaphthene	14.657	153	12011	0.411	ng/ul	99
12) Fluorene	15.647	166	13124	0.398	ng/ul	98
14) Pentachlorophenol	16.978	266	3752	0.693	ng/ul	99
15) Phenanthrene	17.392	178	19476	0.410	ng/ul	99
16) Anthracene	17.485	178	17003	0.415	ng/ul	99
19) Fluoranthene	19.403	202	19632	0.455	ng/ul	99
20) Pyrene	19.766	202	19797	0.467	ng/ul	99
21) Benzo(a)anthracene	21.504	228	12418	0.429	ng/ul	99
22) Chrysene	21.559	228	12965	0.434	ng/ul	100
24) Benzo(b)fluoranthene	23.187	252	12746	0.431	ng/ul	96
25) Benzo(k)fluoranthene	23.237	252	12357	0.430	ng/ul	95
26) Benzo(a)pyrene	23.822	252	10849	0.447	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.424	276	15051	0.451	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.441	278	10697	0.444	ng/ul	99
29) Benzo(g,h,i)perylene	27.199	276	12799	0.456	ng/ul	98

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

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