

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062024\
 Data File : BM046139.D
 Acq On : 21 Jun 2024 00:53
 Operator : MA/JU
 Sample : PB161391BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS391

Quant Time: Jun 21 02:50:08 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM061824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 18 13:58:29 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.463	152	2284	0.400	ng/ul	0.00
4) Naphthalene-d8	10.248	136	7245	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.098	164	4252	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.853	188	8250	0.400	ng/ul	0.00
17) Chrysene-d12	21.032	240	6760	0.400	ng/ul	-0.01
23) Perylene-d12	23.125	264	9117	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.037	96	1842	0.588	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.870	152	3025	0.311	ng/ul	0.00
18) Fluoranthene-d10	18.880	212	6295	0.311	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.067	88	4857	1.382	ng/ul#	77
5) Naphthalene	10.292	128	6429	0.318	ng/ul	98
7) 2-Methylnaphthalene	11.936	142	3543	0.312	ng/ul	99
8) 1-Methylnaphthalene	12.139	142	3752	0.301	ng/ul	99
10) Acenaphthylene	13.825	152	6536	0.324	ng/ul	98
11) Acenaphthene	14.158	153	4581	0.315	ng/ul	97
12) Fluorene	15.172	166	4834	0.309	ng/ul	98
14) Pentachlorophenol	16.540	266	1519	0.614	ng/ul	98
15) Phenanthrene	16.895	178	7232	0.312	ng/ul	99
16) Anthracene	16.996	178	5279	0.306	ng/ul	99
19) Fluoranthene	18.908	202	9895	0.338	ng/ul	99
20) Pyrene	19.266	202	10370	0.328	ng/ul	99
21) Benzo(a)anthracene	21.018	228	8554	0.328	ng/ul	99
22) Chrysene	21.067	228	10921	0.328	ng/ul	99
24) Benzo(b)fluoranthene	22.505	252	10794	0.329	ng/ul	99
25) Benzo(k)fluoranthene	22.546	252	12363	0.317	ng/ul	99
26) Benzo(a)pyrene	23.034	252	10894	0.358	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.182	276	14459	0.290	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.189	278	11301	0.305	ng/ul	99
29) Benzo(g,h,i)perylene	25.796	276	12364	0.289	ng/ul	99

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

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