

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100423\
 Data File : BM042191.D
 Acq On : 04 Oct 2023 16:53
 Operator : MA/JU
 Sample : PB155910BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS910

Quant Time: Oct 05 01:54:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 02 19:28:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.929	152	3407	0.400	ng/ul	0.00
4) Naphthalene-d8	10.746	136	8726	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.583	164	3849	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.337	188	8315	0.400	ng/ul	0.00
17) Chrysene-d12	21.513	240	5453	0.400	ng/ul	0.00
23) Perylene-d12	23.915	264	5547	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.360	96	2422	0.534	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.330	152	2977	0.249	ng/ul	0.00
18) Fluoranthene-d10	19.357	212	5259	0.257	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.393	88	6283	1.326	ng/ul#	79
5) Naphthalene	10.796	128	7473	0.257	ng/ul	98
7) 2-Methylnaphthalene	12.407	142	4291	0.246	ng/ul	96
8) 1-Methylnaphthalene	12.627	142	4301	0.245	ng/ul	97
10) Acenaphthylene	14.315	152	6003	0.247	ng/ul	98
11) Acenaphthene	14.648	153	4802	0.261	ng/ul	99
12) Fluorene	15.638	166	5438	0.262	ng/ul	96
14) Pentachlorophenol	16.974	266	1995	0.529	ng/ul	99
15) Phenanthrene	17.379	178	9248	0.270	ng/ul	99
16) Anthracene	17.472	178	7939	0.253	ng/ul	98
19) Fluoranthene	19.390	202	11497	0.278	ng/ul	97
20) Pyrene	19.752	202	12272	0.279	ng/ul	94
21) Benzo(a)anthracene	21.495	228	9701	0.276	ng/ul	99
22) Chrysene	21.548	228	10543	0.308	ng/ul	99
24) Benzo(b)fluoranthene	23.173	252	11031	0.337	ng/ul	95
25) Benzo(k)fluoranthene	23.222	252	10886	0.338	ng/ul	94
26) Benzo(a)pyrene	23.810	252	9043	0.321	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	26.414	276	12908	0.323	ng/ul#	90
28) Dibenzo(a,h)anthracene	26.431	278	9458	0.314	ng/ul#	92
29) Benzo(g,h,i)perylene	27.186	276	11144	0.334	ng/ul	95

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

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