

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100323\
 Data File : BM042156.D
 Acq On : 03 Oct 2023 10:19
 Operator : MA/JU
 Sample : PB155907BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS907

Quant Time: Oct 03 11:17:49 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	7.937	152	4109	0.400	ng/ul	0.00
4) Naphthalene-d8	10.746	136	9689	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.587	164	4117	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.341	188	8288	0.400	ng/ul	0.00
17) Chrysene-d12	21.515	240	5233	0.400	ng/ul	0.00
23) Perylene-d12	23.921	264	5851	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	3398	0.621	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.330	152	4266	0.321	ng/ul	0.00
18) Fluoranthene-d10	19.361	212	7001	0.356	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.397	88	10333	1.808	ng/ul#	77
5) Naphthalene	10.801	128	10290	0.318	ng/ul	100
7) 2-Methylnaphthalene	12.407	142	6110	0.316	ng/ul	99
8) 1-Methylnaphthalene	12.627	142	5840	0.300	ng/ul	99
10) Acenaphthylene	14.314	152	9189	0.353	ng/ul	99
11) Acenaphthene	14.647	153	6663	0.338	ng/ul	98
12) Fluorene	15.637	166	7338	0.331	ng/ul	98
14) Pentachlorophenol	16.982	266	2827	0.752	ng/ul	99
15) Phenanthrene	17.383	178	11831	0.346	ng/ul	99
16) Anthracene	17.476	178	10364	0.332	ng/ul	99
19) Fluoranthene	19.394	202	14524	0.365	ng/ul	99
20) Pyrene	19.757	202	15415	0.365	ng/ul	95
21) Benzo(a)anthracene	21.498	228	12182	0.361	ng/ul	100
22) Chrysene	21.553	228	12218	0.372	ng/ul	99
24) Benzo(b)fluoranthene	23.175	252	12258	0.355	ng/ul	97
25) Benzo(k)fluoranthene	23.228	252	12854	0.379	ng/ul	95
26) Benzo(a)pyrene	23.813	252	11726	0.394	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.414	276	15886	0.377	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.431	278	11876	0.373	ng/ul	96
29) Benzo(g,h,i)perylene	27.189	276	13337	0.379	ng/ul	97

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

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