

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072624\
 Data File : BN033022.D
 Acq On : 27 Jul 2024 06:52
 Operator : MA/JU
 Sample : PB162203BS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS203

Quant Time: Jul 28 22:56:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN072524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 25 15:43:00 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.545	152	2190	0.400	ng/ul	0.02
4) Naphthalene-d8	10.311	136	9006	0.400	ng/ul	0.01
9) Acenaphthene-d10	14.166	164	5983	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.941	188	12745	0.400	ng/ul	0.00
17) Chrysene-d12	21.154	240	11533	0.400	ng/ul	0.00
23) Perylene-d12	23.343	264	13260	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.064	96	2228	0.804	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.928	152	5061	0.344	ng/ul	0.02
18) Fluoranthene-d10	18.981	212	13399	0.410	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.094	88	6606	2.092	ng/ul#	68
5) Naphthalene	10.366	128	8956	0.387	ng/ul	99
7) 2-Methylnaphthalene	12.000	142	5638	0.360	ng/ul	99
8) 1-Methylnaphthalene	12.198	142	6041	0.357	ng/ul	100
10) Acenaphthylene	13.898	152	9191	0.375	ng/ul	100
11) Acenaphthene	14.236	153	7655	0.416	ng/ul	99
12) Fluorene	15.249	166	8147	0.380	ng/ul	99
14) Pentachlorophenol	16.633	266	1425	0.549	ng/ul	99
15) Phenanthrene	16.983	178	12461	0.365	ng/ul	99
16) Anthracene	17.089	178	11005	0.358	ng/ul	99
19) Fluoranthene	19.009	202	16470	0.434	ng/ul	99
20) Pyrene	19.381	202	17450	0.429	ng/ul	99
21) Benzo(a)anthracene	21.142	228	14486	0.368	ng/ul	99
22) Chrysene	21.192	228	19328	0.421	ng/ul	98
24) Benzo(b)fluoranthene	22.688	252	16155	0.417	ng/ul	94
25) Benzo(k)fluoranthene	22.732	252	20121	0.433	ng/ul#	92
26) Benzo(a)pyrene	23.258	252	16549	0.454	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	25.554	276	19542	0.353	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.554	278	15201	0.346	ng/ul	96
29) Benzo(g,h,i)perylene	26.231	276	15524	0.352	ng/ul	97

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

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