

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102422\
 Data File : BN022307.D
 Acq On : 24 Oct 2022 20:51
 Operator : CG/JU
 Sample : PB148415BS
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS415

Quant Time: Oct 25 04:43:03 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 22 03:40:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.944	152	3898	0.400	ng/ul	0.00
4) Naphthalene-d8	10.772	136	12945	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.619	164	6250	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.374	188	10631	0.400	ng/ul	0.00
17) Chrysene-d12	21.576	240	7625	0.400	ng/ul	0.00
23) Perylene-d12	24.037	264	6827	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.257	96	3417	0.675	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.373	152	6905	0.353	ng/ul	0.00
18) Fluoranthene-d10	19.407	212	9215	0.360	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.291	88	8407	1.616	ng/ul	89
5) Naphthalene	10.822	128	13737	0.359	ng/ul	100
7) 2-Methylnaphthalene	12.444	142	8153	0.338	ng/ul	99
8) 1-Methylnaphthalene	12.664	142	8705	0.351	ng/ul	100
10) Acenaphthylene	14.341	152	9953	0.355	ng/ul	100
11) Acenaphthene	14.679	153	8159	0.345	ng/ul	99
12) Fluorene	15.674	166	8412	0.306	ng/ul	99
14) Pentachlorophenol	17.028	266	1415	0.560	ng/ul	100
15) Phenanthrene	17.417	178	12270	0.344	ng/ul	99
16) Anthracene	17.510	178	10239	0.340	ng/ul	99
19) Fluoranthene	19.440	202	12090	0.346	ng/ul	100
20) Pyrene	19.802	202	12281	0.350	ng/ul	98
21) Benzo(a)anthracene	21.559	228	10539	0.369	ng/ul	99
22) Chrysene	21.611	228	11153	0.368	ng/ul	99
24) Benzo(b)fluoranthene	23.280	252	11607	0.339	ng/ul	95
25) Benzo(k)fluoranthene	23.330	252	11431	0.343	ng/ul	97
26) Benzo(a)pyrene	23.926	252	10241	0.375	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.622	276	13559	0.399	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.642	278	10650	0.392	ng/ul	97
29) Benzo(g,h,i)perylene	27.413	276	11201	0.405	ng/ul	99

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

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