

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062022\
 Data File : BN020295.D
 Acq On : 21 Jun 2022 01:42
 Operator : CG/JU
 Sample : PB145550BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS550

Quant Time: Jun 21 02:39:14 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:13:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.800	152	1126	0.400	ng/ul	0.00
4) Naphthalene-d8	10.584	136	4388	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.423	164	3628	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.162	188	9335	0.400	ng/ul	0.00
17) Chrysene-d12	21.349	240	11634	0.400	ng/ul	0.00
23) Perylene-d12	23.661	264	12497	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	844	0.656	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.174	152	2668	0.368	ng/ul	0.00
18) Fluoranthene-d10	19.192	212	10918	0.455	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.290	88	2129	1.720	ng/ul#	70
5) Naphthalene	10.634	128	4361	0.311	ng/ul	99
7) 2-Methylnaphthalene	12.251	142	3058	0.337	ng/ul	100
8) 1-Methylnaphthalene	12.465	142	3311	0.371	ng/ul	98
10) Acenaphthylene	14.141	152	5618	0.324	ng/ul	99
11) Acenaphthene	14.483	153	4492	0.271	ng/ul	99
12) Fluorene	15.473	166	5553	0.323	ng/ul	99
14) Pentachlorophenol	16.833	266	1269	0.708	ng/ul	99
15) Phenanthrene	17.204	178	10540	0.312	ng/ul	99
16) Anthracene	17.297	178	9042	0.316	ng/ul	98
19) Fluoranthene	19.220	202	14136	0.301	ng/ul	98
20) Pyrene	19.583	202	14959	0.294	ng/ul	98
21) Benzo(a)anthracene	21.334	228	14266	0.316	ng/ul	100
22) Chrysene	21.384	228	16476	0.293	ng/ul	100
24) Benzo(b)fluoranthene	22.959	252	18694	0.210	ng/ul	99
25) Benzo(k)fluoranthene	23.003	252	18772	0.226	ng/ul	99
26) Benzo(a)pyrene	23.556	252	18325	0.324	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.029	276	23285	0.253	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.045	278	18782	0.256	ng/ul	100
29) Benzo(g,h,i)perylene	26.756	276	20616	0.254	ng/ul	99

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

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