

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062022\
 Data File : BN020296.D
 Acq On : 21 Jun 2022 02:18
 Operator : CG/JU
 Sample : PB145553BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS553

Quant Time: Jun 21 03:56:30 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	788	0.400	ng/ul	0.00
4) Naphthalene-d8	10.584	136	3221	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.423	164	2722	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.162	188	7203	0.400	ng/ul	0.00
17) Chrysene-d12	21.349	240	9122	0.400	ng/ul	0.00
23) Perylene-d12	23.661	264	9790	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	689	0.766	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.173	152	2245	0.421	ng/ul	0.00
18) Fluoranthene-d10	19.192	212	9608	0.510	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.290	88	1741	2.010	ng/ul#	70
5) Naphthalene	10.634	128	3582	0.348	ng/ul	98
7) 2-Methylnaphthalene	12.250	142	2566	0.385	ng/ul	99
8) 1-Methylnaphthalene	12.465	142	2813	0.429	ng/ul	100
10) Acenaphthylene	14.141	152	4803	0.369	ng/ul	99
11) Acenaphthene	14.483	153	3831	0.308	ng/ul	98
12) Fluorene	15.469	166	4805	0.373	ng/ul	98
14) Pentachlorophenol	16.833	266	1162	0.840	ng/ul	99
15) Phenanthrene	17.204	178	9180	0.352	ng/ul	99
16) Anthracene	17.297	178	7910	0.359	ng/ul	99
19) Fluoranthene	19.220	202	12509	0.340	ng/ul	98
20) Pyrene	19.583	202	13346	0.334	ng/ul	98
21) Benzo(a)anthracene	21.334	228	12789	0.361	ng/ul	100
22) Chrysene	21.387	228	14662	0.333	ng/ul	99
24) Benzo(b)fluoranthene	22.959	252	16651	0.238	ng/ul	99
25) Benzo(k)fluoranthene	23.003	252	16694	0.256	ng/ul	99
26) Benzo(a)pyrene	23.558	252	16277	0.367	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.032	276	20872	0.290	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.045	278	16879	0.294	ng/ul	99
29) Benzo(g,h,i)perylene	26.750	276	18713	0.294	ng/ul	99

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

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