

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020402.D
 Acq On : 24 Jun 2022 17:51
 Operator : CG/JU
 Sample : PB145731BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS731

Quant Time: Jun 25 00:55:21 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:17:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.754	152	2767	0.400	ng/u1	0.00
4) Naphthalene-d8	10.533	136	9155	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.380	164	6157	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.119	188	14223	0.400	ng/u1	0.00
17) Chrysene-d12	21.312	240	12752	0.400	ng/u1	0.00
23) Perylene-d12	23.595	264	10439	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	2363	0.737	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.128	152	5786	0.385	ng/u1	0.00
18) Fluoranthene-d10	19.149	212	15507	0.369	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.239	88	5890	1.799	ng/u1#	83
5) Naphthalene	10.583	128	9875	0.349	ng/u1	99
7) 2-Methylnaphthalene	12.205	142	6609	0.349	ng/u1	99
8) 1-Methylnaphthalene	12.420	142	6978	0.389	ng/u1	100
10) Acenaphthylene	14.098	152	9925	0.340	ng/u1	99
11) Acenaphthene	14.440	153	8010	0.343	ng/u1	99
12) Fluorene	15.430	166	9547	0.345	ng/u1	99
14) Pentachlorophenol	16.785	266	1446	0.504	ng/u1	99
15) Phenanthrene	17.161	178	17003	0.344	ng/u1	100
16) Anthracene	17.254	178	14637	0.343	ng/u1	100
19) Fluoranthene	19.181	202	19566	0.336	ng/u1	99
20) Pyrene	19.544	202	20199	0.334	ng/u1	98
21) Benzo(a)anthracene	21.294	228	16510	0.360	ng/u1	100
22) Chrysene	21.347	228	18834	0.369	ng/u1	100
24) Benzo(b)fluoranthene	22.905	252	17709	0.367	ng/u1	99
25) Benzo(k)fluoranthene	22.949	252	17482	0.358	ng/u1	100
26) Benzo(a)pyrene	23.495	252	15947	0.363	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	25.926	276	16388	0.375	ng/u1	98
28) Dibenzo(a,h)anthracene	25.946	278	13241	0.375	ng/u1	99
29) Benzo(g,h,i)perylene	26.644	276	13988	0.373	ng/u1	99

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

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