

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062122\
 Data File : BN020332.D
 Acq On : 22 Jun 2022 06:28
 Operator : CG/JU
 Sample : PB145593BS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS593

Quant Time: Jun 22 07:17:38 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	528	0.400	ng/ul	0.00
4) Naphthalene-d8	10.584	136	1907	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.419	164	1515	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.162	188	4346	0.400	ng/ul	0.00
17) Chrysene-d12	21.350	240	7742	0.400	ng/ul	0.00
23) Perylene-d12	23.656	264	9991	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	410	0.680	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.174	152	1200	0.381	ng/ul	0.00
18) Fluoranthene-d10	19.192	212	6175	0.386	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.290	88	1049	1.808	ng/ul#	72
5) Naphthalene	10.628	128	2044	0.335	ng/ul#	95
7) 2-Methylnaphthalene	12.251	142	1338	0.339	ng/ul	96
8) 1-Methylnaphthalene	12.465	142	1470	0.379	ng/ul	99
10) Acenaphthylene	14.141	152	2623	0.362	ng/ul	96
11) Acenaphthene	14.483	153	2004	0.289	ng/ul	99
12) Fluorene	15.469	166	2509	0.350	ng/ul	99
14) Pentachlorophenol	16.829	266	937	1.123	ng/ul	99
15) Phenanthrene	17.204	178	5033	0.320	ng/ul	97
16) Anthracene	17.297	178	4493	0.338	ng/ul	98
19) Fluoranthene	19.220	202	7772	0.249	ng/ul	98
20) Pyrene	19.583	202	8686	0.256	ng/ul	98
21) Benzo(a)anthracene	21.335	228	9968	0.332	ng/ul	98
22) Chrysene	21.385	228	11202	0.299	ng/ul	100
24) Benzo(b)fluoranthene	22.960	252	14238	0.200	ng/ul	97
25) Benzo(k)fluoranthene	23.004	252	14022	0.211	ng/ul	97
26) Benzo(a)pyrene	23.557	252	14682	0.325	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.027	276	19094	0.260	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.044	278	15326	0.261	ng/ul	98
29) Benzo(g,h,i)perylene	26.748	276	17107	0.264	ng/ul	99

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

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