

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020458.D
 Acq On : 26 Jun 2022 05:34
 Operator : CG/JU
 Sample : PB145682BS
 Misc :
 ALS Vial : 76 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS682

Quant Time: Jun 27 03:44:25 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	7.750	152	2506	0.400	ng/ul	0.00
4) Naphthalene-d8	10.528	136	8726	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.376	164	5764	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.119	188	12728	0.400	ng/ul	0.00
17) Chrysene-d12	21.307	240	11905	0.400	ng/ul	0.00
23) Perylene-d12	23.587	264	9934	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.201	96	2049	0.706	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.123	152	5308	0.371	ng/ul	0.00
18) Fluoranthene-d10	19.149	212	13999	0.357	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.235	88	5308	1.790	ng/ul#	84
5) Naphthalene	10.578	128	9363	0.348	ng/ul	100
7) 2-Methylnaphthalene	12.200	142	6296	0.349	ng/ul	99
8) 1-Methylnaphthalene	12.414	142	6680	0.391	ng/ul	99
10) Acenaphthylene	14.093	152	9881	0.362	ng/ul	100
11) Acenaphthene	14.440	153	7552	0.345	ng/ul	98
12) Fluorene	15.426	166	8904	0.344	ng/ul	99
14) Pentachlorophenol	16.785	266	616	0.240	ng/ul	98
15) Phenanthrene	17.161	178	15464	0.349	ng/ul	99
16) Anthracene	17.250	178	13816	0.362	ng/ul	100
19) Fluoranthene	19.177	202	18230	0.335	ng/ul	100
20) Pyrene	19.540	202	19132	0.339	ng/ul	99
21) Benzo(a)anthracene	21.292	228	16886	0.395	ng/ul	99
22) Chrysene	21.342	228	18024	0.379	ng/ul	99
24) Benzo(b)fluoranthene	22.899	252	16982	0.369	ng/ul	98
25) Benzo(k)fluoranthene	22.943	252	16700	0.360	ng/ul	99
26) Benzo(a)pyrene	23.487	252	15788	0.378	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.923	276	15826	0.381	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.937	278	12908	0.384	ng/ul	100
29) Benzo(g,h,i)perylene	26.638	276	13634	0.382	ng/ul	98

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

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