

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092823\
 Data File : BN027980.D
 Acq On : 30 Sep 2023 03:19
 Operator : MA/JU
 Sample : PB155869BS
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS869

Quant Time: Sep 30 03:50:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN091523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 27 03:52:19 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	7242	0.400	ng/ul	0.00
4) Naphthalene-d8	10.779	136	27092	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.601	164	16022	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.354	188	33563	0.400	ng/ul	0.00
17) Chrysene-d12	21.539	240	23878	0.400	ng/ul	0.00
23) Perylene-d12	24.012	264	23541	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.330	96	3921	0.430	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.357	152	11838	0.288	ng/ul	0.00
18) Fluoranthene-d10	19.380	212	23722	0.325	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.363	88	11071	1.166	ng/ul#	84
5) Naphthalene	10.828	128	18221	0.235	ng/ul	100
7) 2-Methylnaphthalene	12.434	142	11927	0.243	ng/ul	100
8) 1-Methylnaphthalene	12.649	142	11588	0.229	ng/ul	99
10) Acenaphthylene	14.328	152	18602	0.266	ng/ul	99
11) Acenaphthene	14.661	153	13194	0.218	ng/ul	97
12) Fluorene	15.651	166	14528	0.204	ng/ul	99
14) Pentachlorophenol	16.991	266	3330	0.424	ng/ul	99
15) Phenanthrene	17.396	178	22713	0.215	ng/ul	99
16) Anthracene	17.489	178	21351	0.242	ng/ul	99
19) Fluoranthene	19.408	202	25060	0.260	ng/ul	97
20) Pyrene	19.775	202	25478	0.256	ng/ul	99
21) Benzo(a)anthracene	21.521	228	21432	0.232	ng/ul	99
22) Chrysene	21.577	228	20169	0.211	ng/ul	99
24) Benzo(b)fluoranthene	23.243	252	19093	0.182	ng/ul	91
25) Benzo(k)fluoranthene	23.296	252	19680	0.193	ng/ul#	89
26) Benzo(a)pyrene	23.898	252	17916	0.212	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.619	276	22140	0.205	ng/ul#	92
28) Dibenzo(a,h)anthracene	26.636	278	18182	0.209	ng/ul	95
29) Benzo(g,h,i)perylene	27.434	276	18147	0.198	ng/ul	96

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

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