

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092623\
 Data File : BN027873.D
 Acq On : 26 Sep 2023 22:09
 Operator : MA/JU
 Sample : PB155617BS
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS617

Quant Time: Sep 27 03:02:05 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 : Fri Sep 22 03:58:00 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.966	152	2641	0.400	ng/ul	0.00
4) Naphthalene-d8	10.784	136	8973	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.606	164	6165	0.400	ng/ul	0.01
13) Phenanthrene-d10	17.358	188	15064	0.400	ng/ul	0.02
17) Chrysene-d12	21.539	240	15114	0.400	ng/ul	0.02
23) Perylene-d12	24.012	264	16539	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	2608	0.784	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.363	152	6074	0.447	ng/ul	0.00
18) Fluoranthene-d10	19.380	212	17649	0.382	ng/ul	0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.368	88	7595	2.194	ng/ul#	85
5) Naphthalene	10.834	128	9613	0.375	ng/ul	99
7) 2-Methylnaphthalene	12.440	142	6551	0.403	ng/ul	100
8) 1-Methylnaphthalene	12.660	142	6462	0.385	ng/ul	98
10) Acenaphthylene	14.333	152	10155	0.377	ng/ul	100
11) Acenaphthene	14.670	153	8128	0.349	ng/ul	97
12) Fluorene	15.651	166	9764	0.357	ng/ul	99
14) Pentachlorophenol	16.987	266	2913	0.826	ng/ul	99
15) Phenanthrene	17.401	178	16716	0.353	ng/ul	98
16) Anthracene	17.494	178	14556	0.367	ng/ul	98
19) Fluoranthene	19.413	202	20197	0.330	ng/ul	100
20) Pyrene	19.780	202	21211	0.337	ng/ul	98
21) Benzo(a)anthracene	21.522	228	21607	0.369	ng/ul	99
22) Chrysene	21.577	228	21910	0.363	ng/ul	99
24) Benzo(b)fluoranthene	23.240	252	23808	0.323	ng/ul	97
25) Benzo(k)fluoranthene	23.296	252	24149	0.338	ng/ul	98
26) Benzo(a)pyrene	23.898	252	22327	0.377	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.606	276	27919	0.367	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.626	278	22779	0.372	ng/ul	98
29) Benzo(g,h,i)perylene	27.418	276	23418	0.363	ng/ul	98

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

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