

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092623\
 Data File : BN027861.D
 Acq On : 26 Sep 2023 14:18
 Operator : MA/JU
 Sample : PB155730BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS730

Quant Time: Sep 27 03:00:26 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	7.967	152	2809	0.400	ng/ul	0.00
4) Naphthalene-d8	10.784	136	9693	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.610	164	6499	0.400	ng/ul	0.02
13) Phenanthrene-d10	17.359	188	15516	0.400	ng/ul	0.02
17) Chrysene-d12	21.545	240	16030	0.400	ng/ul	0.02
23) Perylene-d12	24.021	264	18568	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	3021	0.854	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.368	152	7337	0.500	ng/ul	0.01
18) Fluoranthene-d10	19.385	212	20887	0.427	ng/ul	0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.368	88	8887	2.413	ng/ul#	86
5) Naphthalene	10.839	128	11332	0.409	ng/ul	99
7) 2-Methylnaphthalene	12.440	142	7716	0.439	ng/ul	100
8) 1-Methylnaphthalene	12.660	142	7598	0.419	ng/ul	99
10) Acenaphthylene	14.337	152	12333	0.434	ng/ul	100
11) Acenaphthene	14.671	153	9441	0.385	ng/ul	99
12) Fluorene	15.656	166	11212	0.389	ng/ul	99
14) Pentachlorophenol	16.991	266	2846	0.784	ng/ul	98
15) Phenanthrene	17.401	178	19196	0.394	ng/ul	99
16) Anthracene	17.494	178	17289	0.424	ng/ul	99
19) Fluoranthene	19.413	202	23469	0.362	ng/ul	98
20) Pyrene	19.780	202	24644	0.369	ng/ul	99
21) Benzo(a)anthracene	21.530	228	25500	0.411	ng/ul	99
22) Chrysene	21.583	228	24785	0.387	ng/ul	99
24) Benzo(b)fluoranthene	23.252	252	27388	0.331	ng/ul	96
25) Benzo(k)fluoranthene	23.302	252	27682	0.345	ng/ul	96
26) Benzo(a)pyrene	23.907	252	26253	0.394	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.619	276	33960	0.398	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.640	278	27851	0.406	ng/ul	98
29) Benzo(g,h,i)perylene	27.431	276	28059	0.387	ng/ul	98

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

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