

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092223\
 Data File : BN027820.D
 Acq On : 22 Sep 2023 11:46
 Operator : MA/JU
 Sample : 04332-19
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Sep 22 12:26:07 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	8.004	152	5415	0.400	ng/ul	0.00
4) Naphthalene-d8	10.823	136	18772	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.638	164	11612	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.392	188	26185	0.400	ng/ul	0.00
17) Chrysene-d12	21.574	240	22200	0.400	ng/ul	0.00
23) Perylene-d12	24.067	264	19178	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.380	96	16532	2.424	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.401	152	4416	0.155	ng/ul	0.00
18) Fluoranthene-d10	19.413	212	11088	0.164	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.414	88	156	0.022	ng/ul#	1
5) Naphthalene	10.872	128	2184	0.041	ng/ul#	85
7) 2-Methylnaphthalene	12.473	142	1679	0.049	ng/ul	100
8) 1-Methylnaphthalene	12.693	142	1488	0.042	ng/ul	98
10) Acenaphthylene	14.370	152	1414	0.028	ng/ul#	74
15) Phenanthrene	17.434	178	12603	0.153	ng/ul	98
16) Anthracene	17.527	178	1679	0.024	ng/ul#	74
19) Fluoranthene	19.445	202	32853	0.366	ng/ul	100
20) Pyrene	19.812	202	28909	0.313	ng/ul	98
21) Benzo(a)anthracene	21.553	228	12657	0.147	ng/ul	94
22) Chrysene	21.609	228	15341	0.173	ng/ul#	93
24) Benzo(b)fluoranthene	23.295	252	27967	0.327	ng/ul#	24
26) Benzo(a)pyrene	23.950	252	12175	0.177	ng/ul#	61
27) Indeno(1,2,3-cd)pyrene	26.683	276	8660	0.098	ng/ul#	88
28) Dibenzo(a,h)anthracene	26.706	278	2082	0.029	ng/ul#	1
29) Benzo(g,h,i)perylene	27.518	276	8151	0.109	ng/ul#	62

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

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