

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092223\
 Data File : BN027822.D
 Acq On : 22 Sep 2023 12:59
 Operator : MA/JU
 Sample : 04332-18
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AE5

Quant Time: Sep 22 13:56:16 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	6140	0.400	ng/ul	0.00
4) Naphthalene-d8	10.823	136	21141	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.638	164	12858	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.392	188	29436	0.400	ng/ul	0.00
17) Chrysene-d12	21.574	240	23752	0.400	ng/ul	# 0.00
23) Perylene-d12	24.068	264	20870	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.380	96	17755	2.296	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.401	152	5246	0.164	ng/ul	0.00
18) Fluoranthene-d10	19.413	212	13707	0.189	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.872	128	1990	0.033	ng/ul#	83
7) 2-Methylnaphthalene	12.478	142	1737	0.045	ng/ul	99
8) 1-Methylnaphthalene	12.693	142	1473	0.037	ng/ul	99
10) Acenaphthylene	14.370	152	1779	0.032	ng/ul#	79
15) Phenanthrene	17.434	178	13482	0.146	ng/ul	97
16) Anthracene	17.523	178	1975	0.026	ng/ul#	69
19) Fluoranthene	19.446	202	35017	0.365	ng/ul	96
20) Pyrene	19.813	202	30853	0.312	ng/ul	99
21) Benzo(a)anthracene	21.557	228	13760	0.150	ng/ul#	91
22) Chrysene	21.612	228	17306	0.182	ng/ul#	92
24) Benzo(b)fluoranthene	23.293	252	32485	0.349	ng/ul#	1
26) Benzo(a)pyrene	23.954	252	12590	0.168	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.694	276	10370	0.108	ng/ul#	1
28) Dibenzo(a,h)anthracene	26.707	278	2735	0.035	ng/ul#	1
29) Benzo(g,h,i)perylene	27.522	276	11517	0.141	ng/ul#	34

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

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