

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN103124\
 Data File : BN034765.D
 Acq On : 31 Oct 2024 14:40
 Operator : RC/JU
 Sample : P4476-02DL2 20X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 XA101DL2

Quant Time: Oct 31 15:05:49 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 31 11:13:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.604	152	9271	0.400	ng/ul	0.00
4) Naphthalene-d8	10.366	136	28441	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.235	164	15825	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.983	188	32299	0.400	ng/ul	0.00
17) Chrysene-d12	21.174	240	22852	0.400	ng/ul	0.00
23) Perylene-d12	23.358	264	20857	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.178	96	279	0.029	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.967	152	594	0.017	ng/ul	0.00
18) Fluoranthene-d10	19.014	212	1512	0.020	ng/ul	0.00
Target Compounds						Qvalue
5) Naphthalene	10.416	128	31909	0.424	ng/ul	100
10) Acenaphthylene	13.953	152	11312	0.157	ng/ul	99
11) Acenaphthene	14.300	153	23326	0.423	ng/ul	99
12) Fluorene	15.290	166	11966	0.195	ng/ul	99
15) Phenanthrene	17.021	178	16118	0.165	ng/ul	99
16) Anthracene	17.114	178	71672	0.793	ng/ul	100
19) Fluoranthene	19.042	202	25161	0.227	ng/ul	98
20) Pyrene	19.404	202	12154	0.105	ng/ul	96
21) Benzo(a)anthracene	21.160	228	21427	0.234	ng/ul	100
22) Chrysene	21.209	228	8431	0.092	ng/ul	98
24) Benzo(b)fluoranthene	22.708	252	17538	0.207	ng/ul	95
25) Benzo(k)fluoranthene	22.752	252	21181	0.247	ng/ul	97
26) Benzo(a)pyrene	23.264	252	5720	0.078	ng/ul#	69
27) Indeno(1,2,3-cd)pyrene	25.533	276	29207	0.375	ng/ul	98
28) Dibenzo(a,h)anthracene	25.550	278	20752	0.353	ng/ul	99
29) Benzo(g,h,i)perylene	26.191	276	17768	0.291	ng/ul	99

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

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