

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080423\
 Data File : BN026818.D
 Acq On : 04 Aug 2023 14:52
 Operator : MA/JU
 Sample : 03805-10DL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X3H69DL

Quant Time: Aug 04 22:30:53 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 04 03:46:36 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.131	152	8596	0.400	ng/ul	0.01
4) Naphthalene-d8	10.955	136	25463	0.400	ng/ul	0.01
9) Acenaphthene-d10	14.754	164	15417	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.498	188	34534	0.400	ng/ul	0.00
17) Chrysene-d12	21.674	240	30065	0.400	ng/ul	0.00
23) Perylene-d12	24.231	264	32420	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.460	96	447	0.044	ng/ul	0.02
6) 2-Methylnaphthalene-d10	12.528	152	1105	0.029	ng/ul	0.01
18) Fluoranthene-d10	19.511	212	3570	0.038	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	11.010	128	105775	1.470	ng/ul	99
10) Acenaphthylene	14.481	152	119132	1.379	ng/ul	99
11) Acenaphthene	14.814	153	70812	1.185	ng/ul	99
12) Fluorene	15.795	166	44523	0.658	ng/ul	99
15) Phenanthrene	17.540	178	78126	0.689	ng/ul	100
16) Anthracene	17.629	178	55650	0.543	ng/ul	99
19) Fluoranthene	19.543	202	106728	0.855	ng/ul	100
20) Pyrene	19.906	202	130492	0.979	ng/ul	98
21) Benzo(a)anthracene	21.656	228	35088	0.293	ng/ul	100
22) Chrysene	21.712	228	63143	0.535	ng/ul	99
24) Benzo(b)fluoranthene	23.439	252	74785	0.610	ng/ul	98
25) Benzo(k)fluoranthene	23.489	252	63169	0.522	ng/ul	98
26) Benzo(a)pyrene	24.114	252	85207	0.770	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.942	276	92782	0.641	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.975	278	59342	0.514	ng/ul	98
29) Benzo(g,h,i)perylene	27.780	276	62641	0.520	ng/ul	99

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

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