

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111924\
 Data File : BN035195.D
 Acq On : 20 Nov 2024 12:43
 Operator : RC/JU
 Sample : P4875-14MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E2A01MS

Quant Time: Nov 20 13:29:26 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 18 00:44:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.541	152	3417	0.400	ng/ul	0.00
4) Naphthalene-d8	10.301	136	8109	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.180	164	5074	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.937	188	10364	0.400	ng/ul	0.00
17) Chrysene-d12	21.148	240	9033	0.400	ng/ul	0.00
23) Perylene-d12	23.320	264	9886	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	638	0.204	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.901	152	3149	0.265	ng/ul	-0.01
18) Fluoranthene-d10	18.977	212	9464	0.380	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.170	88	1889	0.557	ng/ul	98
5) Naphthalene	10.350	128	5019	0.246	ng/ul	99
7) 2-Methylnaphthalene	11.972	142	3387	0.239	ng/ul	99
8) 1-Methylnaphthalene	12.198	142	3467	0.242	ng/ul	99
10) Acenaphthylene	13.898	152	5225	0.260	ng/ul	99
11) Acenaphthene	14.245	153	3996	0.259	ng/ul	98
12) Fluorene	15.240	166	4904	0.271	ng/ul	100
14) Pentachlorophenol	16.591	266	1861	0.771	ng/ul	98
15) Phenanthrene	16.975	178	8634	0.315	ng/ul	99
16) Anthracene	17.068	178	7438	0.292	ng/ul	99
19) Fluoranthene	19.005	202	10939	0.348	ng/ul	99
20) Pyrene	19.367	202	11460	0.350	ng/ul	98
21) Benzo(a)anthracene	21.131	228	11088	0.354	ng/ul	99
22) Chrysene	21.183	228	11490	0.364	ng/ul	99
24) Benzo(b)fluoranthene	22.674	252	11763	0.360	ng/ul	100
25) Benzo(k)fluoranthene	22.715	252	12251	0.364	ng/ul	99
26) Benzo(a)pyrene	23.226	252	9890	0.331	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.473	276	14316	0.355	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.490	278	11252	0.356	ng/ul	99
29) Benzo(g,h,i)perylene	26.131	276	11264	0.354	ng/ul	99

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

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