

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN103124\
 Data File : BN034771.D
 Acq On : 31 Oct 2024 18:16
 Operator : RC/JU
 Sample : P4435-02MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHK68MS

Quant Time: Nov 01 05:02:51 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	6275	0.400	ng/ul	0.00
4) Naphthalene-d8	10.372	136	19479	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.236	164	10728	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.983	188	21882	0.400	ng/ul	0.00
17) Chrysene-d12	21.172	240	15486	0.400	ng/ul	0.00
23) Perylene-d12	23.355	264	12883	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.174	96	2849	0.439	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.967	152	6980	0.286	ng/ul	0.00
18) Fluoranthene-d10	19.014	212	16386	0.325	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.208	88	7865	1.103	ng/ul#	91
5) Naphthalene	10.416	128	13939	0.270	ng/ul	99
7) 2-Methylnaphthalene	12.038	142	9252	0.276	ng/ul	99
8) 1-Methylnaphthalene	12.264	142	10993	0.322	ng/ul	100
10) Acenaphthylene	13.954	152	15908	0.326	ng/ul	99
11) Acenaphthene	14.301	153	10867	0.291	ng/ul	99
12) Fluorene	15.291	166	12660	0.304	ng/ul	100
14) Pentachlorophenol	16.633	266	847	0.196	ng/ul	99
15) Phenanthrene	17.021	178	33705	0.510	ng/ul	99
16) Anthracene	17.110	178	19485	0.318	ng/ul	99
19) Fluoranthene	19.042	202	45016	0.599	ng/ul	98
20) Pyrene	19.405	202	47501	0.607	ng/ul	97
21) Benzo(a)anthracene	21.158	228	28066	0.453	ng/ul	98
22) Chrysene	21.210	228	29073	0.468	ng/ul	99
24) Benzo(b)fluoranthene	22.706	252	26216	0.500	ng/ul	99
25) Benzo(k)fluoranthene	22.750	252	21754	0.411	ng/ul	97
26) Benzo(a)pyrene	23.262	252	20707	0.459	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.531	276	18674	0.388	ng/ul	97
28) Dibenzo(a,h)anthracene	25.548	278	12730	0.350	ng/ul	99
29) Benzo(g,h,i)perylene	26.192	276	15471	0.410	ng/ul	99

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

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