

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034968.D
 Acq On : 10 Nov 2024 07:59
 Operator : RC/JU
 Sample : P4746-10MSD
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GCP83MSD

Quant Time: Nov 10 21:51:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 09 00:09:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.567	152	2713	0.400	ng/ul	0.00
4) Naphthalene-d8	10.328	136	6533	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.203	164	4762	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.950	188	11385	0.400	ng/ul	0.00
17) Chrysene-d12	21.143	240	11812	0.400	ng/ul	0.00
23) Perylene-d12	23.305	264	12749	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	620	0.240	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.928	152	3411	0.341	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	12834	0.423	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.183	88	2114	0.747	ng/ul#	93
5) Naphthalene	10.378	128	5751	0.352	ng/ul	99
7) 2-Methylnaphthalene	12.000	142	4144	0.340	ng/ul	99
8) 1-Methylnaphthalene	12.225	142	4179	0.341	ng/ul	99
10) Acenaphthylene	13.917	152	6817	0.350	ng/ul	99
11) Acenaphthene	14.264	153	5194	0.352	ng/ul	99
12) Fluorene	15.254	166	6548	0.349	ng/ul	99
14) Pentachlorophenol	16.603	266	2329	0.885	ng/ul	98
15) Phenanthrene	16.988	178	11795	0.388	ng/ul	99
16) Anthracene	17.081	178	10655	0.373	ng/ul	99
19) Fluoranthene	19.014	202	15505	0.393	ng/ul	98
20) Pyrene	19.372	202	16404	0.401	ng/ul	97
21) Benzo(a)anthracene	21.125	228	16331	0.382	ng/ul	100
22) Chrysene	21.178	228	16655	0.382	ng/ul	100
24) Benzo(b)fluoranthene	22.662	252	17219	0.384	ng/ul	99
25) Benzo(k)fluoranthene	22.703	252	17910	0.382	ng/ul	99
26) Benzo(a)pyrene	23.212	252	15368	0.384	ng/ul#	99
27) Indeno(1,2,3-cd)pyrene	25.453	276	20950	0.428	ng/ul	94
28) Dibenzo(a,h)anthracene	25.470	278	16306	0.427	ng/ul	97
29) Benzo(g,h,i)perylene	26.107	276	16253	0.434	ng/ul	98

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

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