

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070823\
 Data File : BN026397.D
 Acq On : 08 Jul 2023 22:17
 Operator : MA/JU
 Sample : 03350-04MSD
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YBW27MSD

Quant Time: Jul 08 23:54:31 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 08 09:20:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.930	152	4687	0.400	ng/ul	0.00
4) Naphthalene-d8	10.737	136	16355	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.570	164	9256	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.326	188	15827	0.400	ng/ul	0.00
17) Chrysene-d12	21.517	240	9831	0.400	ng/ul	-0.01
23) Perylene-d12	23.958	264	9875	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	2455	0.465	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.332	152	5870	0.233	ng/ul	0.00
18) Fluoranthene-d10	19.354	212	10114	0.270	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.335	88	6188	1.104	ng/ul#	87
5) Naphthalene	10.787	128	12225	0.261	ng/ul	98
7) 2-Methylnaphthalene	12.403	142	7194	0.241	ng/ul	100
8) 1-Methylnaphthalene	12.618	142	7714	0.249	ng/ul	99
10) Acenaphthylene	14.293	152	11719	0.270	ng/ul	99
11) Acenaphthene	14.635	153	9223	0.260	ng/ul	99
12) Fluorene	15.625	166	9648	0.237	ng/ul	99
14) Pentachlorophenol	16.984	266	1112	0.196	ng/ul	99
15) Phenanthrene	17.364	178	13615	0.260	ng/ul	97
16) Anthracene	17.461	178	13222	0.285	ng/ul	96
19) Fluoranthene	19.382	202	14252	0.271	ng/ul	100
20) Pyrene	19.744	202	15238	0.279	ng/ul	99
21) Benzo(a)anthracene	21.500	228	10425	0.257	ng/ul	99
22) Chrysene	21.552	228	11365	0.269	ng/ul	97
24) Benzo(b)fluoranthene	23.212	252	11111	0.264	ng/ul	84
25) Benzo(k)fluoranthene	23.259	252	11547	0.273	ng/ul#	87
26) Benzo(a)pyrene	23.853	252	9933	0.271	ng/ul#	82
27) Indeno(1,2,3-cd)pyrene	26.514	276	12218	0.306	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.524	278	9294	0.307	ng/ul	91
29) Benzo(g,h,i)perylene	27.295	276	10905	0.307	ng/ul	97

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

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