

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071122\
 Data File : BN020704.D
 Acq On : 11 Jul 2022 16:02
 Operator : CG/JU
 Sample : N3577-04MSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YB324MSD

Quant Time: Jul 11 18:39:14 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 02 00:47:40 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.825	152	4625	0.400	ng/ul	0.00
4) Naphthalene-d8	10.623	136	14680	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.474	164	8038	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.225	188	15507	0.400	ng/ul	0.00
17) Chrysene-d12	21.431	240	11469	0.400	ng/ul	-0.01
23) Perylene-d12	23.784	264	7984	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.227	96	1079	0.201	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.223	152	6298	0.261	ng/ul	0.00
18) Fluoranthene-d10	19.262	212	15119	0.400	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.260	88	3245	0.593	ng/ul	91
5) Naphthalene	10.672	128	12703	0.280	ng/ul	99
7) 2-Methylnaphthalene	12.294	142	7556	0.249	ng/ul	99
8) 1-Methylnaphthalene	12.514	142	7997	0.278	ng/ul	100
10) Acenaphthylene	14.192	152	10397	0.273	ng/ul	99
11) Acenaphthene	14.534	153	8474	0.278	ng/ul	99
12) Fluorene	15.529	166	9433	0.261	ng/ul	99
14) Pentachlorophenol	16.888	266	1490	0.477	ng/ul	98
15) Phenanthrene	17.268	178	16434	0.305	ng/ul	100
16) Anthracene	17.361	178	13025	0.280	ng/ul	100
19) Fluoranthene	19.295	202	19960	0.381	ng/ul	100
20) Pyrene	19.657	202	20364	0.375	ng/ul	99
21) Benzo(a)anthracene	21.416	228	13552	0.329	ng/ul	100
22) Chrysene	21.469	228	16037	0.350	ng/ul	99
24) Benzo(b)fluoranthene	23.070	252	12917	0.350	ng/ul	99
25) Benzo(k)fluoranthene	23.117	252	13737	0.368	ng/ul	99
26) Benzo(a)pyrene	23.678	252	11311	0.337	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	26.216	276	11875	0.355	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.237	278	9646	0.357	ng/ul	99
29) Benzo(g,h,i)perylene	26.961	276	10884	0.380	ng/ul	99

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

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