

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062122\
 Data File : BN020330.D
 Acq On : 22 Jun 2022 05:16
 Operator : CG/JU
 Sample : N3359-08MS
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EW4R7MS

Quant Time: Jun 22 05:45:19 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:13:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.792	152	897	0.400	ng/ul	0.00
4) Naphthalene-d8	10.579	136	4141	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.419	164	2666	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.158	188	6425	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.347	240	9355	0.400	ng/ul	# 0.00
23) Perylene-d12	23.656	264	10983	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	246	0.240	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.168	152	1926	0.281	ng/ul	-0.01
18) Fluoranthene-d10	19.188	212	9154	0.474	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.290	88	729	0.739	ng/ul#	29
5) Naphthalene	10.628	128	3551	0.268	ng/ul	99
7) 2-Methylnaphthalene	12.245	142	2332	0.272	ng/ul	100
8) 1-Methylnaphthalene	12.460	142	2464	0.292	ng/ul	97
10) Acenaphthylene	14.136	152	4467	0.350	ng/ul	99
11) Acenaphthene	14.479	153	3311	0.272	ng/ul	99
12) Fluorene	15.464	166	3905	0.309	ng/ul	97
14) Pentachlorophenol	16.820	266	1849	1.498	ng/ul	100
15) Phenanthrene	17.196	178	7706	0.331	ng/ul	96
16) Anthracene	17.289	178	8378	0.426	ng/ul	98
19) Fluoranthene	19.216	202	11262	0.298	ng/ul	95
20) Pyrene	19.583	202	12361	0.302	ng/ul	98
21) Benzo(a)anthracene	21.330	228	13199	0.363	ng/ul	96
22) Chrysene	21.385	228	13512	0.299	ng/ul	98
24) Benzo(b)fluoranthene	22.955	252	16330	0.208	ng/ul	94
25) Benzo(k)fluoranthene	23.001	252	15555	0.213	ng/ul	94
26) Benzo(a)pyrene	23.554	252	16198	0.326	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.024	276	20621	0.255	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.037	278	16757	0.260	ng/ul	98
29) Benzo(g,h,i)perylene	26.745	276	18248	0.256	ng/ul	99

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

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