

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062122\
 Data File : BN020331.D
 Acq On : 22 Jun 2022 05:52
 Operator : CG/JU
 Sample : N3359-09MSD
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EW4R7MSD

Quant Time: Jun 22 06:36:25 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	7.796	152	991	0.400	ng/ul	0.00
4) Naphthalene-d8	10.577	136	4540	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.417	164	2914	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.157	188	6913	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.350	240	9853	0.400	ng/ul	0.00
23) Perylene-d12	23.656	264	11419	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	249	0.220	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.167	152	2061	0.275	ng/ul	-0.01
18) Fluoranthene-d10	19.186	212	9801	0.482	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.290	88	808	0.742	ng/ul#	37
5) Naphthalene	10.627	128	3962	0.273	ng/ul	99
7) 2-Methylnaphthalene	12.244	142	2498	0.266	ng/ul	99
8) 1-Methylnaphthalene	12.458	142	2652	0.287	ng/ul	98
10) Acenaphthylene	14.135	152	4815	0.345	ng/ul	99
11) Acenaphthene	14.477	153	3630	0.272	ng/ul	99
12) Fluorene	15.463	166	4181	0.303	ng/ul	99
14) Pentachlorophenol	16.823	266	2046	1.541	ng/ul	97
15) Phenanthrene	17.199	178	8297	0.332	ng/ul	97
16) Anthracene	17.287	178	8928	0.422	ng/ul	98
19) Fluoranthene	19.219	202	11996	0.302	ng/ul	97
20) Pyrene	19.581	202	12969	0.301	ng/ul	99
21) Benzo(a)anthracene	21.333	228	13833	0.362	ng/ul	96
22) Chrysene	21.382	228	14105	0.296	ng/ul	97
24) Benzo(b)fluoranthene	22.957	252	17142	0.210	ng/ul	94
25) Benzo(k)fluoranthene	23.001	252	16331	0.215	ng/ul	95
26) Benzo(a)pyrene	23.554	252	16956	0.328	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.024	276	21445	0.255	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.034	278	17501	0.261	ng/ul	97
29) Benzo(g,h,i)perylene	26.748	276	18997	0.256	ng/ul	99

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

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