

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061223\
 Data File : BN025783.D
 Acq On : 12 Jun 2023 15:56
 Operator : MA/JU
 Sample : 02951-05MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EW999MS

Quant Time: Jun 13 00:02:43 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN060923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 10 02:01:49 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	9921	0.400	ng/ul	0.00
4) Naphthalene-d8	10.820	136	29007	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.640	164	17150	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.386	188	35755	0.400	ng/ul	0.00
17) Chrysene-d12	21.571	240	20934	0.400	ng/ul	0.00
23) Perylene-d12	24.055	264	17133	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.356	96	3067	0.314	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.409	152	7517	0.204	ng/ul	0.00
18) Fluoranthene-d10	19.411	212	16833	0.255	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.390	88	7927	0.759	ng/ul	94
5) Naphthalene	10.875	128	13900	0.188	ng/ul	98
7) 2-Methylnaphthalene	12.480	142	9155	0.206	ng/ul	98
8) 1-Methylnaphthalene	12.695	142	9728	0.199	ng/ul	99
10) Acenaphthylene	14.363	152	12559	0.185	ng/ul	96
11) Acenaphthene	14.705	153	10973	0.194	ng/ul	99
12) Fluorene	15.691	166	12980	0.206	ng/ul	98
14) Pentachlorophenol	17.052	266	149	0.039	ng/ul#	85
15) Phenanthrene	17.428	178	25254	0.242	ng/ul	98
16) Anthracene	17.521	178	17860	0.203	ng/ul	96
19) Fluoranthene	19.439	202	28469	0.297	ng/ul	98
20) Pyrene	19.801	202	30631	0.309	ng/ul	98
21) Benzo(a)anthracene	21.553	228	17742	0.266	ng/ul	95
22) Chrysene	21.606	228	20514	0.269	ng/ul	96
24) Benzo(b)fluoranthene	23.292	252	15877	0.227	ng/ul	80
25) Benzo(k)fluoranthene	23.339	252	15768	0.224	ng/ul#	81
26) Benzo(a)pyrene	23.944	252	12371	0.208	ng/ul#	77
27) Indeno(1,2,3-cd)pyrene	26.659	276	14902	0.220	ng/ul#	44
28) Dibenzo(a,h)anthracene	26.679	278	11760	0.229	ng/ul	92
29) Benzo(g,h,i)perylene	27.457	276	14029	0.231	ng/ul	96

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

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