

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062724\
 Data File : BN032274.D
 Acq On : 27 Jun 2024 10:30
 Operator : MA/JU
 Sample : SSTD0.202
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.2231

Quant Time: Jun 27 12:49:38 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN062724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 27 12:48:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.620	152	7151	0.400	ng/ul	0.00
4) Naphthalene-d8	10.402	136	21410	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.266	164	14197	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.027	188	33406	0.400	ng/ul	0.00
17) Chrysene-d12	21.228	240	28238	0.400	ng/ul	0.00
23) Perylene-d12	23.449	264	29033	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.156	96	1646	0.217	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.003	152	6355	0.189	ng/ul	0.00
18) Fluoranthene-d10	19.063	212	17382	0.207	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.190	88	2029	0.238	ng/ul#	87
5) Naphthalene	10.452	128	11315	0.203	ng/ul	98
7) 2-Methylnaphthalene	12.080	142	7390	0.192	ng/ul	100
8) 1-Methylnaphthalene	12.294	142	7595	0.191	ng/ul	98
10) Acenaphthylene	13.988	152	11748	0.199	ng/ul	99
11) Acenaphthene	14.330	153	9236	0.208	ng/ul	98
12) Fluorene	15.330	166	10955	0.203	ng/ul	99
14) Pentachlorophenol	16.711	266	1356	0.177	ng/ul	98
15) Phenanthrene	17.065	178	18194	0.199	ng/ul	99
16) Anthracene	17.162	178	16087	0.196	ng/ul	97
19) Fluoranthene	19.090	202	21745	0.211	ng/ul	98
20) Pyrene	19.458	202	22941	0.213	ng/ul	99
21) Benzo(a)anthracene	21.213	228	20328	0.194	ng/ul	99
22) Chrysene	21.266	228	22897	0.216	ng/ul	99
24) Benzo(b)fluoranthene	22.782	252	25993	0.210	ng/ul	92
25) Benzo(k)fluoranthene	22.826	252	25853	0.216	ng/ul#	93
26) Benzo(a)pyrene	23.355	252	17670	0.201	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	25.698	276	29540	0.243	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.708	278	21867	0.234	ng/ul	95
29) Benzo(g,h,i)perylene	26.383	276	23587	0.243	ng/ul	99

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

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