

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061822\
 Data File : BN020267.D
 Acq On : 18 Jun 2022 20:41
 Operator : CG/JU
 Sample : SSTD1.606
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.6206

Quant Time: Jun 20 00:58:03 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 00:54:39 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.796	152	954	0.400	ng/ul	0.00
4) Naphthalene-d8	10.584	136	3553	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.423	164	2655	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.162	188	6634	0.400	ng/ul	0.00
17) Chrysene-d12	21.353	240	9283	0.400	ng/ul	0.00
23) Perylene-d12	23.662	264	8364	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	1666	1.532	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.173	152	9776	1.724	ng/ul	0.00
18) Fluoranthene-d10	19.192	212	34042	1.213	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.290	88	1762	1.541	ng/ul#	74
5) Naphthalene	10.628	128	15304	1.369	ng/ul	94
7) 2-Methylnaphthalene	12.245	142	10900	1.517	ng/ul	98
8) 1-Methylnaphthalene	12.465	142	11850	1.768	ng/ul	100
10) Acenaphthylene	14.141	152	17783	1.350	ng/ul	97
11) Acenaphthene	14.483	153	14179	1.400	ng/ul	98
12) Fluorene	15.469	166	17654	1.516	ng/ul	99
14) Pentachlorophenol	16.829	266	1770	1.269	ng/ul	99
15) Phenanthrene	17.204	178	31895	1.355	ng/ul	98
16) Anthracene	17.297	178	28066	1.379	ng/ul	97
19) Fluoranthene	19.220	202	43142	1.088	ng/ul	99
20) Pyrene	19.583	202	44153	1.106	ng/ul	98
21) Benzo(a)anthracene	21.335	228	43473	1.278	ng/ul	99
22) Chrysene	21.388	228	49039	1.277	ng/ul	98
24) Benzo(b)fluoranthene	22.963	252	55591	1.448	ng/ul	94
25) Benzo(k)fluoranthene	23.007	252	54629	1.449	ng/ul	95
26) Benzo(a)pyrene	23.563	252	49196	1.400	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	26.034	276	67391	1.683	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.050	278	55180	1.688	ng/ul	97
29) Benzo(g,h,i)perylene	26.751	276	59003	1.741	ng/ul	99

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

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