

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052422\
 Data File : BN019948.D
 Acq On : 25 May 2022 13:09
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4365

Quant Time: May 26 01:15:17 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN052322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 23 14:58:28 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	4404	0.400	ng/ul	0.00
4) Naphthalene-d8	10.639	136	15007	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.470	164	8775	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.213	188	17200	0.400	ng/ul	0.00
17) Chrysene-d12	21.397	240	11061	0.400	ng/ul	0.00
23) Perylene-d12	23.735	264	9023	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	1656	0.403	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.229	152	9662	0.394	ng/ul	0.00
18) Fluoranthene-d10	19.239	212	17120	0.433	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.353	88	1682	0.393	ng/ul#	87
5) Naphthalene	10.689	128	17048	0.362	ng/ul	99
7) 2-Methylnaphthalene	12.300	142	11283	0.362	ng/ul#	100
8) 1-Methylnaphthalene	12.520	142	11440	0.394	ng/ul#	100
10) Acenaphthylene	14.192	152	17167	0.343	ng/ul	99
11) Acenaphthene	14.534	153	12236	0.361	ng/ul	99
12) Fluorene	15.520	166	13630	0.360	ng/ul	99
14) Pentachlorophenol	16.867	266	944	0.330	ng/ul	89
15) Phenanthrene	17.251	178	21852	0.361	ng/ul	99
16) Anthracene	17.344	178	19530	0.344	ng/ul	99
19) Fluoranthene	19.267	202	22287	0.404	ng/ul	100
20) Pyrene	19.629	202	21829	0.390	ng/ul	100
21) Benzo(a)anthracene	21.382	228	15607	0.345	ng/ul	100
22) Chrysene	21.435	228	16462	0.360	ng/ul	100
24) Benzo(b)fluoranthene	23.028	252	14870	0.373	ng/ul	92
25) Benzo(k)fluoranthene	23.075	252	14324	0.369	ng/ul	97
26) Benzo(a)pyrene	23.633	252	14072	0.373	ng/ul#	78
27) Indeno(1,2,3-cd)pyrene	26.138	276	15229	0.385	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.155	278	12350	0.381	ng/ul#	80
29) Benzo(g,h,i)perylene	26.872	276	12982	0.384	ng/ul	98

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

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