

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071624\
 Data File : BN032790.D
 Acq On : 17 Jul 2024 22:04
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4265

Quant Time: Jul 17 23:33:04 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 17 02:35:25 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.548	152	3564	0.400	ng/ul	0.00
4) Naphthalene-d8	10.316	136	11230	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.193	164	6003	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.961	188	11078	0.400	ng/ul	0.00
17) Chrysene-d12	21.171	240	10575	0.400	ng/ul	# 0.00
23) Perylene-d12	23.368	264	13066	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.101	96	1797	0.420	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.922	152	6072	0.358	ng/ul	-0.01
18) Fluoranthene-d10	18.999	212	11108	0.322	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.131	88	1838	0.377	ng/ul#	69
5) Naphthalene	10.365	128	11180	0.371	ng/ul	100
7) 2-Methylnaphthalene	11.993	142	6687	0.355	ng/ul	98
8) 1-Methylnaphthalene	12.213	142	7163	0.368	ng/ul	100
10) Acenaphthylene	13.920	152	10213	0.421	ng/ul	99
11) Acenaphthene	14.258	153	7243	0.361	ng/ul	98
12) Fluorene	15.262	166	7446	0.325	ng/ul	99
14) Pentachlorophenol	16.640	266	863	0.444	ng/ul	93
15) Phenanthrene	17.003	178	10650	0.353	ng/ul	99
16) Anthracene	17.100	178	10143	0.412	ng/ul	99
19) Fluoranthene	19.031	202	13614	0.317	ng/ul	99
20) Pyrene	19.399	202	14297	0.321	ng/ul	99
21) Benzo(a)anthracene	21.156	228	13578	0.403	ng/ul	96
22) Chrysene	21.209	228	15275	0.314	ng/ul	96
24) Benzo(b)fluoranthene	22.711	252	15096	0.327	ng/ul	80
25) Benzo(k)fluoranthene	22.751	252	15820	0.282	ng/ul#	79
26) Benzo(a)pyrene	23.275	252	14938	0.374	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	25.586	276	19011	0.352	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.586	278	15381	0.367	ng/ul#	84
29) Benzo(g,h,i)perylene	26.263	276	15588	0.345	ng/ul#	90

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

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