

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111523\
 Data File : BN028644.D
 Acq On : 16 Nov 2023 01:33
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4313

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/17/2023
 Supervised By :mohammad ahmed 11/20/2023

Quant Time: Nov 16 13:54:39 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 16 07:25:53 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.772	152	5456m	0.400	ng/ul	0.09
4) Naphthalene-d8	10.497	136	16589m	0.400	ng/ul	0.03
9) Acenaphthene-d10	14.327	164	11262	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.091	188	15119	0.400	ng/ul	0.01
17) Chrysene-d12	21.296	240	18813	0.400	ng/ul	0.00
23) Perylene-d12	23.581	264	19460	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	3134m	0.520	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.109	152	7211m	0.289	ng/ul	0.04
18) Fluoranthene-d10	19.119	212	23302	0.385	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.156	88	3017m	0.453	ng/ul	
5) Naphthalene	10.541	128	15918	0.342	ng/ul	96
7) 2-Methylnaphthalene	12.180	142	9173m	0.299	ng/ul	
8) 1-Methylnaphthalene	12.373	142	11383	0.366	ng/ul	100
10) Acenaphthylene	14.054	152	21766	0.387	ng/ul	98
11) Acenaphthene	14.392	153	17908	0.433	ng/ul	99
12) Fluorene	15.396	166	14937	0.310	ng/ul	98
14) Pentachlorophenol	16.758	266	1411	0.389	ng/ul	99
15) Phenanthrene	17.134	178	21851	0.470	ng/ul	98
16) Anthracene	17.239	178	11624	0.273	ng/ul	95
19) Fluoranthene	19.152	202	31474	0.393	ng/ul	100
20) Pyrene	19.514	202	36644	0.442	ng/ul	98
21) Benzo(a)anthracene	21.278	228	22097	0.297	ng/ul	98
22) Chrysene	21.331	228	28394	0.390	ng/ul	99
24) Benzo(b)fluoranthene	22.888	252	28071	0.384	ng/ul	92
25) Benzo(k)fluoranthene	22.935	252	34234	0.448	ng/ul	94
26) Benzo(a)pyrene	23.485	252	29048	0.421	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	25.927	276	38327	0.501	ng/ul#	45
28) Dibenzo(a,h)anthracene	25.954	278	28663	0.469	ng/ul	96
29) Benzo(g,h,i)perylene	26.642	276	34061	0.543	ng/ul	99

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

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