

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111524\
 Data File : BN035097.D
 Acq On : 15 Nov 2024 09:49
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4337

Quant Time: Nov 15 11:51:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 00:41:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.549	152	2739	0.400	ng/ul	0.00
4) Naphthalene-d8	10.311	136	6726	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.189	164	5146	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.937	188	12956	0.400	ng/ul	0.00
17) Chrysene-d12	21.136	240	11545	0.400	ng/ul	# 0.00
23) Perylene-d12	23.299	264	11303	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	1046	0.401	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.912	152	4391	0.426	ng/ul	0.00
18) Fluoranthene-d10	18.972	212	14161	0.478	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.174	88	1219	0.427	ng/ul	96
5) Naphthalene	10.361	128	7126	0.423	ng/ul	98
7) 2-Methylnaphthalene	11.983	142	5283	0.421	ng/ul	100
8) 1-Methylnaphthalene	12.203	142	5339	0.423	ng/ul	98
10) Acenaphthylene	13.902	152	8484	0.404	ng/ul	100
11) Acenaphthene	14.249	153	6500	0.408	ng/ul	99
12) Fluorene	15.244	166	8284	0.409	ng/ul	99
14) Pentachlorophenol	16.591	266	1396	0.466	ng/ul	98
15) Phenanthrene	16.979	178	14076	0.407	ng/ul	99
16) Anthracene	17.068	178	13369	0.411	ng/ul	98
19) Fluoranthene	19.000	202	17839	0.462	ng/ul#	95
20) Pyrene	19.362	202	18965	0.474	ng/ul#	94
21) Benzo(a)anthracene	21.119	228	17192	0.411	ng/ul	100
22) Chrysene	21.172	228	16591	0.389	ng/ul	99
24) Benzo(b)fluoranthene	22.656	252	15658	0.394	ng/ul	97
25) Benzo(k)fluoranthene	22.700	252	16063	0.386	ng/ul#	93
26) Benzo(a)pyrene	23.206	252	14557	0.410	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	25.450	276	17092	0.394	ng/ul#	91
28) Dibenzo(a,h)anthracene	25.463	278	13333	0.394	ng/ul#	97
29) Benzo(g,h,i)perylene	26.100	276	13621	0.411	ng/ul#	97

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

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