

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111524\
 Data File : BN035132.D
 Acq On : 16 Nov 2024 08:02
 Operator : RC/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4339

Quant Time: Nov 18 00:22:19 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	7.550	152	2848	0.400	ng/ul	0.00
4) Naphthalene-d8	10.311	136	7496	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.185	164	6025	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.933	188	14318	0.400	ng/ul	0.00
17) Chrysene-d12	21.134	240	10778	0.400	ng/ul	# 0.00
23) Perylene-d12	23.296	264	9576	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	1079	0.398	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.906	152	4906	0.427	ng/ul	-0.01
18) Fluoranthene-d10	18.967	212	14204	0.513	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.174	88	1203	0.405	ng/ul	96
5) Naphthalene	10.361	128	7878	0.420	ng/ul	98
7) 2-Methylnaphthalene	11.983	142	5864	0.419	ng/ul	98
8) 1-Methylnaphthalene	12.203	142	6022	0.428	ng/ul	100
10) Acenaphthylene	13.898	152	9876	0.401	ng/ul	99
11) Acenaphthene	14.245	153	7683	0.412	ng/ul	97
12) Fluorene	15.240	166	9440	0.398	ng/ul	100
14) Pentachlorophenol	16.591	266	886	0.268	ng/ul	98
15) Phenanthrene	16.975	178	15732	0.411	ng/ul	99
16) Anthracene	17.068	178	14478	0.403	ng/ul	98
19) Fluoranthene	19.000	202	18074	0.502	ng/ul#	96
20) Pyrene	19.362	202	18369	0.492	ng/ul#	95
21) Benzo(a)anthracene	21.116	228	15582	0.399	ng/ul	100
22) Chrysene	21.169	228	15621	0.392	ng/ul	99
24) Benzo(b)fluoranthene	22.653	252	13724	0.408	ng/ul	96
25) Benzo(k)fluoranthene	22.694	252	14090	0.400	ng/ul#	96
26) Benzo(a)pyrene	23.200	252	12406	0.412	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.443	276	15192	0.413	ng/ul#	90
28) Dibenzo(a,h)anthracene	25.457	278	11860	0.414	ng/ul#	96
29) Benzo(g,h,i)perylene	26.094	276	11887	0.423	ng/ul#	96

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

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