

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121324\
 Data File : BN035611.D
 Acq On : 14 Dec 2024 00:18
 Operator : RC/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4367

Quant Time: Dec 14 00:49:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 10 16:10:57 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.289	152	2830	0.400	ng/ul	0.00
4) Naphthalene-d8	10.037	136	8419	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.949	164	5575	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.717	188	10868	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.964	240	7792	0.400	ng/ul	# 0.00
23) Perylene-d12	23.054	264	8949	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.959	96	1028	0.397	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.637	152	5043	0.409	ng/ul	0.00
18) Fluoranthene-d10	18.772	212	10403	0.484	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.993	88	1038	0.370	ng/ul#	68
5) Naphthalene	10.086	128	8621	0.406	ng/ul	99
7) 2-Methylnaphthalene	11.714	142	5869	0.399	ng/ul	98
8) 1-Methylnaphthalene	11.939	142	5994	0.404	ng/ul	98
10) Acenaphthylene	13.662	152	9110	0.412	ng/ul	99
11) Acenaphthene	14.014	153	6811	0.403	ng/ul	98
12) Fluorene	15.018	166	7531	0.379	ng/ul	99
14) Pentachlorophenol	16.388	266	893	0.353	ng/ul	97
15) Phenanthrene	16.760	178	11417	0.397	ng/ul	99
16) Anthracene	16.852	178	10771	0.404	ng/ul	99
19) Fluoranthene	18.800	202	12731	0.469	ng/ul#	94
20) Pyrene	19.167	202	13709	0.486	ng/ul#	91
21) Benzo(a)anthracene	20.947	228	11137	0.412	ng/ul	96
22) Chrysene	20.999	228	11311	0.416	ng/ul	97
24) Benzo(b)fluoranthene	22.443	252	11052	0.374	ng/ul#	81
25) Benzo(k)fluoranthene	22.481	252	12208	0.400	ng/ul#	81
26) Benzo(a)pyrene	22.963	252	11043	0.408	ng/ul#	77
27) Indeno(1,2,3-cd)pyrene	25.084	276	14010	0.384	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.104	278	10866	0.380	ng/ul#	81
29) Benzo(g,h,i)perylene	25.698	276	11255	0.391	ng/ul#	83

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

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