

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111124\
 Data File : BN035009.D
 Acq On : 11 Nov 2024 17:37
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4332

Quant Time: Nov 11 18:06:19 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 : Sat Nov 09 00:09:12 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.562	152	3200	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.328	136	8160	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.199	164	5253	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	10964	0.400	ng/ul	0.00
17) Chrysene-d12	21.143	240	10254	0.400	ng/ul	# 0.00
23) Perylene-d12	23.308	264	11428	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	1007	0.331	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	4776	0.382	ng/ul	-0.01
18) Fluoranthene-d10	18.977	212	10802	0.410	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.183	88	1138	0.341	ng/ul	97
5) Naphthalene	10.372	128	7994	0.391	ng/ul	98
7) 2-Methylnaphthalene	12.000	142	5608	0.368	ng/ul	99
8) 1-Methylnaphthalene	12.220	142	5637	0.368	ng/ul	100
10) Acenaphthylene	13.912	152	8126	0.379	ng/ul	99
11) Acenaphthene	14.264	153	6210	0.382	ng/ul	99
12) Fluorene	15.254	166	7137	0.345	ng/ul	99
14) Pentachlorophenol	16.599	266	973	0.384	ng/ul	97
15) Phenanthrene	16.988	178	11241	0.384	ng/ul	99
16) Anthracene	17.076	178	10578	0.384	ng/ul	99
19) Fluoranthene	19.009	202	13423	0.392	ng/ul#	96
20) Pyrene	19.372	202	14044	0.395	ng/ul#	95
21) Benzo(a)anthracene	21.125	228	14204	0.382	ng/ul	100
22) Chrysene	21.178	228	13872	0.366	ng/ul	98
24) Benzo(b)fluoranthene	22.662	252	14126	0.352	ng/ul	94
25) Benzo(k)fluoranthene	22.703	252	14959	0.356	ng/ul#	95
26) Benzo(a)pyrene	23.212	252	13635	0.380	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	25.460	276	17235	0.392	ng/ul#	92
28) Dibenzo(a,h)anthracene	25.473	278	13586	0.397	ng/ul#	97
29) Benzo(g,h,i)perylene	26.111	276	13398	0.399	ng/ul#	97

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

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