

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112224\
 Data File : BN035237.D
 Acq On : 22 Nov 2024 11:51
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4347

Quant Time: Nov 22 12:45:36 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 : Fri Nov 22 12:45:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.318	152	3592	0.400	ng/ul	0.00
4) Naphthalene-d8	10.069	136	8950	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.981	164	5738	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.747	188	13715	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.993	240	11064	0.400	ng/ul	# 0.00
23) Perylene-d12	23.100	264	9771	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.976	96	1444	0.440	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.670	152	4924	0.375	ng/ul	0.00
18) Fluoranthene-d10	18.800	212	14210	0.466	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.010	88	1657	0.465	ng/ul#	71
5) Naphthalene	10.119	128	9277	0.411	ng/ul	99
7) 2-Methylnaphthalene	11.747	142	6372	0.407	ng/ul	98
8) 1-Methylnaphthalene	11.972	142	6461	0.409	ng/ul	98
10) Acenaphthylene	13.690	152	10537	0.463	ng/ul	99
11) Acenaphthene	14.041	153	7601	0.436	ng/ul	98
12) Fluorene	15.045	166	9123	0.446	ng/ul	99
14) Pentachlorophenol	16.405	266	1667	0.522	ng/ul	99
15) Phenanthrene	16.789	178	15599	0.430	ng/ul	99
16) Anthracene	16.882	178	14958	0.444	ng/ul	99
19) Fluoranthene	18.833	202	19311	0.501	ng/ul#	93
20) Pyrene	19.200	202	20642	0.515	ng/ul#	91
21) Benzo(a)anthracene	20.979	228	17013	0.443	ng/ul	100
22) Chrysene	21.028	228	16793	0.435	ng/ul	100
24) Benzo(b)fluoranthene	22.481	252	14371	0.445	ng/ul	94
25) Benzo(k)fluoranthene	22.521	252	15253	0.458	ng/ul#	94
26) Benzo(a)pyrene	23.010	252	13061	0.443	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	25.151	276	15604	0.391	ng/ul#	75
28) Dibenzo(a,h)anthracene	25.171	278	11643	0.373	ng/ul#	92
29) Benzo(g,h,i)perylene	25.775	276	11955	0.380	ng/ul#	92

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

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN11624.M
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