

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032724\
 Data File : BN030500.D
 Acq On : 27 Mar 2024 12:04
 Operator : MA/JU
 Sample : SSTD0.281
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.2210

Quant Time: Mar 27 15:07:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN032724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 27 15:02:37 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.724	152	4993	0.400	ng/ul	0.00
4) Naphthalene-d8	10.507	136	15093	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.358	164	8797	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.108	188	20783	0.400	ng/ul	0.00
17) Chrysene-d12	21.301	240	16010	0.400	ng/ul	0.00
23) Perylene-d12	23.549	264	13599	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.218	96	1227	0.224	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.102	152	4508	0.198	ng/ul	0.00
18) Fluoranthene-d10	19.137	212	10016	0.202	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	1139	0.205	ng/ul#	88
5) Naphthalene	10.557	128	8177	0.199	ng/ul	99
7) 2-Methylnaphthalene	12.174	142	5339	0.196	ng/ul	99
8) 1-Methylnaphthalene	12.394	142	5687	0.198	ng/ul	100
10) Acenaphthylene	14.081	152	7743	0.195	ng/ul	98
11) Acenaphthene	14.423	153	6117	0.198	ng/ul	98
12) Fluorene	15.413	166	7156	0.196	ng/ul	98
14) Pentachlorophenol	16.749	266	672	0.177	ng/ul	98
15) Phenanthrene	17.146	178	12328	0.198	ng/ul	99
16) Anthracene	17.239	178	10041	0.191	ng/ul	99
19) Fluoranthene	19.170	202	13155	0.199	ng/ul	100
20) Pyrene	19.532	202	13181	0.203	ng/ul	100
21) Benzo(a)anthracene	21.283	228	10621	0.192	ng/ul	98
22) Chrysene	21.336	228	12284	0.201	ng/ul	99
24) Benzo(b)fluoranthene	22.870	252	10934	0.192	ng/ul	93
25) Benzo(k)fluoranthene	22.914	252	11081	0.185	ng/ul#	94
26) Benzo(a)pyrene	23.446	252	8593	0.186	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	25.823	276	9578	0.183	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.843	278	7472	0.180	ng/ul	95
29) Benzo(g,h,i)perylene	26.517	276	8456	0.190	ng/ul	97

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

