

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071223\
 Data File : BN026481.D
 Acq On : 12 Jul 2023 19:50
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV236

Quant Time: Jul 13 04:35:54 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN071223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 13 04:33:34 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.009	152	2355	0.400	ng/ul	0.00
4) Naphthalene-d8	10.809	136	11219	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.626	164	5537	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.372	188	8207	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.558	240	5798	0.400	ng/ul	0.00
23) Perylene-d12	23.999	264	6145	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.343	96	1736	0.560	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.398	152	5556	0.364	ng/ul	0.00
18) Fluoranthene-d10	19.396	212	6779	0.337	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.381	88	1641	0.559	ng/ul	97
5) Naphthalene	10.858	128	12218	0.397	ng/ul	99
7) 2-Methylnaphthalene	12.469	142	6572	0.376	ng/ul	100
8) 1-Methylnaphthalene	12.678	142	7279	0.389	ng/ul	100
10) Acenaphthylene	14.348	152	9901	0.365	ng/ul	99
11) Acenaphthene	14.686	153	8194	0.397	ng/ul	99
12) Fluorene	15.676	166	7673	0.354	ng/ul	100
14) Pentachlorophenol	17.039	266	692	0.334	ng/ul	99
15) Phenanthrene	17.414	178	10096	0.386	ng/ul	100
16) Anthracene	17.511	178	8457	0.373	ng/ul	100
19) Fluoranthene	19.428	202	8896	0.327	ng/ul	99
20) Pyrene	19.791	202	9846	0.334	ng/ul	98
21) Benzo(a)anthracene	21.540	228	8088	0.380	ng/ul	99
22) Chrysene	21.593	228	9528	0.393	ng/ul	99
24) Benzo(b)fluoranthene	23.250	252	9429	0.394	ng/ul	96
25) Benzo(k)fluoranthene	23.297	252	9756	0.389	ng/ul#	97
26) Benzo(a)pyrene	23.888	252	8688	0.408	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.551	276	11465	0.410	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.564	278	8892	0.415	ng/ul	98
29) Benzo(g,h,i)perylene	27.332	276	10147	0.411	ng/ul	99

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

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