

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020624\
 Data File : BN029536.D
 Acq On : 06 Feb 2024 14:07
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV229

Quant Time: Feb 07 02:18:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 02:17:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	4986	0.400	ng/ul	0.00
4) Naphthalene-d8	10.689	136	13557	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.530	164	5951	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.285	188	11225	0.400	ng/ul	0.00
17) Chrysene-d12	21.484	240	7729	0.400	ng/ul	0.00
23) Perylene-d12	23.861	264	7863	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.336	96	2311	0.372	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.284	152	6236	0.362	ng/ul	0.00
18) Fluoranthene-d10	19.318	212	9376	0.369	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.370	88	2503	0.370	ng/ul	99
5) Naphthalene	10.738	128	14432	0.378	ng/ul	100
7) 2-Methylnaphthalene	12.355	142	8326	0.381	ng/ul	100
8) 1-Methylnaphthalene	12.575	142	7918	0.355	ng/ul	100
10) Acenaphthylene	14.252	152	11159	0.384	ng/ul	100
11) Acenaphthene	14.594	153	8176	0.370	ng/ul	98
12) Fluorene	15.584	166	8739	0.379	ng/ul	98
14) Pentachlorophenol	16.930	266	770	0.308	ng/ul#	49
15) Phenanthrene	17.323	178	13054	0.372	ng/ul	99
16) Anthracene	17.420	178	10898	0.369	ng/ul	99
19) Fluoranthene	19.346	202	13221	0.372	ng/ul	99
20) Pyrene	19.708	202	13642	0.375	ng/ul	99
21) Benzo(a)anthracene	21.467	228	10124	0.363	ng/ul	100
22) Chrysene	21.520	228	11630	0.381	ng/ul	99
24) Benzo(b)fluoranthene	23.139	252	10837	0.366	ng/ul	99
25) Benzo(k)fluoranthene	23.186	252	12352	0.377	ng/ul	98
26) Benzo(a)pyrene	23.756	252	10041	0.361	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.322	276	13004	0.385	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.349	278	9987	0.377	ng/ul	98
29) Benzo(g,h,i)perylene	27.063	276	11431	0.373	ng/ul	99

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

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