

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040723\
 Data File : BN024817.D
 Acq On : 07 Apr 2023 18:44
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV208

Quant Time: Apr 08 00:54:19 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN040723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 08 00:52:14 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	6499	0.400	ng/ul	0.00
4) Naphthalene-d8	10.765	136	19148	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.593	164	10774	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.338	188	23496	0.400	ng/ul	0.00
17) Chrysene-d12	21.529	240	20678	0.400	ng/ul	0.00
23) Perylene-d12	23.973	264	20164	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.318	96	2594	0.366	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.354	152	10563	0.430	ng/ul	0.00
18) Fluoranthene-d10	19.368	212	22815	0.444	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.352	88	2809	0.359	ng/ul	95
5) Naphthalene	10.814	128	19517	0.400	ng/ul	99
7) 2-Methylnaphthalene	12.425	142	12747	0.420	ng/ul	100
8) 1-Methylnaphthalene	12.640	142	12051	0.375	ng/ul	99
10) Acenaphthylene	14.316	152	16908	0.418	ng/ul	99
11) Acenaphthene	14.653	153	13836	0.399	ng/ul	100
12) Fluorene	15.639	166	15710	0.405	ng/ul	99
14) Pentachlorophenol	16.992	266	1553	0.310	ng/ul	98
15) Phenanthrene	17.381	178	25994	0.393	ng/ul	100
16) Anthracene	17.473	178	21520	0.391	ng/ul	100
19) Fluoranthene	19.396	202	28622	0.402	ng/ul	98
20) Pyrene	19.763	202	29643	0.406	ng/ul	98
21) Benzo(a)anthracene	21.511	228	26571	0.407	ng/ul	99
22) Chrysene	21.567	228	28916	0.409	ng/ul	100
24) Benzo(b)fluoranthene	23.224	252	31658	0.391	ng/ul	97
25) Benzo(k)fluoranthene	23.274	252	30273	0.384	ng/ul	98
26) Benzo(a)pyrene	23.864	252	26889	0.401	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.510	276	31546	0.399	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.531	278	24552	0.395	ng/ul	100
29) Benzo(g,h,i)perylene	27.285	276	25984	0.388	ng/ul	99

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

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