

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034919.D
 Acq On : 09 Nov 2024 00:16
 Operator : RC/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV222

Quant Time: Nov 09 00:41:30 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 09 00:09:12 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	2526	0.400	ng/ul	0.00
4) Naphthalene-d8	10.339	136	6407	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.208	164	4845	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.954	188	12162	0.400	ng/ul	0.00
17) Chrysene-d12	21.148	240	12886	0.400	ng/ul	0.00
23) Perylene-d12	23.314	264	13406	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	985	0.410	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.934	152	4222	0.430	ng/ul	0.00
18) Fluoranthene-d10	18.986	212	13842	0.418	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.187	88	1077	0.409	ng/ul	96
5) Naphthalene	10.388	128	6808	0.425	ng/ul	98
7) 2-Methylnaphthalene	12.011	142	5016	0.419	ng/ul	100
8) 1-Methylnaphthalene	12.231	142	4690	0.390	ng/ul	98
10) Acenaphthylene	13.926	152	8418	0.425	ng/ul	99
11) Acenaphthene	14.268	153	5975	0.398	ng/ul	99
12) Fluorene	15.263	166	7691	0.403	ng/ul	99
14) Pentachlorophenol	16.607	266	1085	0.386	ng/ul	98
15) Phenanthrene	16.996	178	13559	0.417	ng/ul	99
16) Anthracene	17.085	178	12627	0.413	ng/ul	98
19) Fluoranthene	19.018	202	17541	0.407	ng/ul	99
20) Pyrene	19.381	202	18611	0.417	ng/ul	99
21) Benzo(a)anthracene	21.134	228	18626	0.399	ng/ul	99
22) Chrysene	21.183	228	19704	0.414	ng/ul	100
24) Benzo(b)fluoranthene	22.671	252	19324	0.410	ng/ul	98
25) Benzo(k)fluoranthene	22.712	252	20471	0.415	ng/ul	98
26) Benzo(a)pyrene	23.217	252	17448	0.414	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.470	276	20941	0.407	ng/ul	97
28) Dibenzo(a,h)anthracene	25.487	278	16391	0.408	ng/ul	99
29) Benzo(g,h,i)perylene	26.124	276	15826	0.402	ng/ul	99

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

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