

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122721\
 Data File : BN018213.D
 Acq On : 27 Dec 2021 21:03
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV214

Quant Time: Dec 28 00:37:36 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 28 00:34:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.607	152	4303	0.400	ng/ul	0.00
4) Naphthalene-d8	10.388	136	14816	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.249	164	10621	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.991	188	24318	0.400	ng/ul	0.00
17) Chrysene-d12	21.183	240	25800	0.400	ng/ul	0.00
23) Perylene-d12	23.404	264	24884	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.072	96	1452	0.361	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.983	152	9702	0.377	ng/ul	0.00
18) Fluoranthene-d10	19.023	212	29993	0.392	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.105	88	1733	0.388	ng/ul#	74
5) Naphthalene	10.432	128	17269	0.393	ng/ul	99
7) 2-Methylnaphthalene	12.060	142	12488	0.388	ng/ul	100
8) 1-Methylnaphthalene	12.280	142	11918	0.368	ng/ul	99
10) Acenaphthylene	13.963	152	19906	0.434	ng/ul	99
11) Acenaphthene	14.310	153	15022	0.406	ng/ul	99
12) Fluorene	15.300	166	17948	0.393	ng/ul	99
14) Pentachlorophenol	16.653	266	3215	0.334	ng/ul	99
15) Phenanthrene	17.033	178	30785	0.383	ng/ul	99
16) Anthracene	17.122	178	26818	0.371	ng/ul	99
19) Fluoranthene	19.055	202	40021	0.390	ng/ul	100
20) Pyrene	19.418	202	41466	0.387	ng/ul	99
21) Benzo(a)anthracene	21.168	228	36975	0.376	ng/ul	100
22) Chrysene	21.221	228	41843	0.392	ng/ul	99
24) Benzo(b)fluoranthene	22.737	252	39473	0.379	ng/ul	100
25) Benzo(k)fluoranthene	22.781	252	46797	0.394	ng/ul	100
26) Benzo(a)pyrene	23.307	252	40857	0.414	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.661	276	46452	0.370	ng/ul	99
28) Dibenzo(a,h)anthracene	25.681	278	36201	0.373	ng/ul	99
29) Benzo(g,h,i)perylene	26.348	276	40133	0.375	ng/ul	100

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

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