

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122121\
 Data File : BN018140.D
 Acq On : 21 Dec 2021 23:35
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV207

Quant Time: Dec 22 14:09:24 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 22 14:08:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.611	152	4524	0.400	ng/ul	0.00
4) Naphthalene-d8	10.388	136	14126	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.249	164	8682	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.995	188	19121	0.400	ng/ul	0.00
17) Chrysene-d12	21.192	240	18285	0.400	ng/ul	0.00
23) Perylene-d12	23.421	264	15450	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.080	96	1800	0.385	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.994	152	7873	0.405	ng/ul	0.00
18) Fluoranthene-d10	19.032	212	21430	0.432	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.114	88	2486	0.440	ng/ul	90
5) Naphthalene	10.438	128	16625	0.414	ng/ul	99
7) 2-Methylnaphthalene	12.066	142	10615	0.411	ng/ul	100
8) 1-Methylnaphthalene	12.286	142	10462	0.390	ng/ul	100
10) Acenaphthylene	13.972	152	15682	0.457	ng/ul	99
11) Acenaphthene	14.314	153	12516	0.414	ng/ul	99
12) Fluorene	15.304	166	14313	0.408	ng/ul	99
14) Pentachlorophenol	16.662	266	1104	0.307	ng/ul	98
15) Phenanthrene	17.038	178	25499	0.408	ng/ul	100
16) Anthracene	17.130	178	20057	0.397	ng/ul	100
19) Fluoranthene	19.060	202	30190	0.430	ng/ul	99
20) Pyrene	19.422	202	31171	0.424	ng/ul	99
21) Benzo(a)anthracene	21.177	228	22996	0.378	ng/ul	100
22) Chrysene	21.227	228	31526	0.423	ng/ul	99
24) Benzo(b)fluoranthene	22.749	252	25986	0.439	ng/ul	100
25) Benzo(k)fluoranthene	22.796	252	31457	0.427	ng/ul	99
26) Benzo(a)pyrene	23.322	252	24683	0.441	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.681	276	25931	0.351	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.697	278	19598	0.351	ng/ul	98
29) Benzo(g,h,i)perylene	26.368	276	22560	0.371	ng/ul	99

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

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