

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN083021\
 Data File : BN016112.D
 Acq On : 30 Aug 2021 15:30
 Operator : CG/JU
 Sample : SST00.108
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.1208

Quant Time: Aug 30 17:21:13 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN083021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 30 17:16:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	1395	0.400	ng/ul	0.00
4) Naphthalene-d8	10.661	136	5882	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.497	164	3459	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.242	188	7486	0.400	ng/ul	0.00
17) Chrysene-d12	21.439	240	7609	0.400	ng/ul	0.00
23) Perylene-d12	23.827	264	7191	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.314	96	231	0.122	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.250	152	953	0.108	ng/ul	0.00
18) Fluoranthene-d10	19.275	212	2159	0.105	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.344	88	262	0.137	ng/ul#	68
5) Naphthalene	10.710	128	1549	0.102	ng/ul#	91
7) 2-Methylnaphthalene	12.321	142	1031	0.100	ng/ul	97
8) 1-Methylnaphthalene	12.541	142	1042	0.104	ng/ul	99
10) Acenaphthylene	14.219	152	1331	0.109	ng/ul#	92
11) Acenaphthene	14.561	153	1087	0.103	ng/ul	97
12) Fluorene	15.547	166	1280	0.103	ng/ul#	95
14) Pentachlorophenol	16.896	266	220	0.140	ng/ul	94
15) Phenanthrene	17.284	178	2264	0.110	ng/ul	94
16) Anthracene	17.377	178	1838	0.103	ng/ul	93
19) Fluoranthene	19.303	202	2441	0.103	ng/ul#	94
20) Pyrene	19.666	202	2473	0.103	ng/ul#	94
21) Benzo(a)anthracene	21.421	228	2512	0.106	ng/ul	95
22) Chrysene	21.471	228	2611	0.102	ng/ul	97
24) Benzo(b)fluoranthene	23.099	252	2667	0.099	ng/ul	79
25) Benzo(k)fluoranthene	23.146	252	2615	0.097	ng/ul#	85
26) Benzo(a)pyrene	23.719	252	2277	0.101	ng/ul#	75
27) Indeno(1,2,3-cd)pyrene	26.303	276	2906	0.097	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.320	278	2402	0.096	ng/ul#	84
29) Benzo(g,h,i)perylene	27.054	276	2499	0.097	ng/ul#	89

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

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