

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090721\
 Data File : BN016232.D
 Acq On : 07 Sep 2021 11:58
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4269

Quant Time: Sep 07 12:58:18 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 07 12:55:06 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	2260	0.400	ng/ul	0.00
4) Naphthalene-d8	10.655	136	9588	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.497	164	5984	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.242	188	12658	0.400	ng/ul	0.00
17) Chrysene-d12	21.430	240	12582	0.400	ng/ul	0.00
23) Perylene-d12	23.815	264	12156	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.302	96	970	0.369	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.250	152	5649	0.418	ng/ul	0.00
18) Fluoranthene-d10	19.271	212	13933	0.413	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.340	88	1011	0.352	ng/ul#	91
5) Naphthalene	10.705	128	9724	0.381	ng/ul	100
7) 2-Methylnaphthalene	12.321	142	6789	0.408	ng/ul	100
8) 1-Methylnaphthalene	12.541	142	6774	0.408	ng/ul	99
10) Acenaphthylene	14.214	152	9390	0.415	ng/ul	99
11) Acenaphthene	14.557	153	7259	0.383	ng/ul	99
12) Fluorene	15.547	166	8537	0.394	ng/ul	99
14) Pentachlorophenol	16.900	266	592	0.289	ng/ul	95
15) Phenanthrene	17.284	178	13916	0.381	ng/ul	99
16) Anthracene	17.377	178	13088	0.415	ng/ul	99
19) Fluoranthene	19.303	202	16547	0.395	ng/ul	99
20) Pyrene	19.666	202	17162	0.404	ng/ul	95
21) Benzo(a)anthracene	21.415	228	16165	0.420	ng/ul	98
22) Chrysene	21.465	228	16171	0.372	ng/ul	98
24) Benzo(b)fluoranthene	23.090	252	16519	0.361	ng/ul	93
25) Benzo(k)fluoranthene	23.137	252	16453	0.358	ng/ul	93
26) Benzo(a)pyrene	23.710	252	14739	0.380	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.290	276	18120	0.361	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.306	278	14991	0.363	ng/ul	94
29) Benzo(g,h,i)perylene	27.048	276	15520	0.357	ng/ul	96

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

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