

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111524\
 Data File : BN035114.D
 Acq On : 15 Nov 2024 20:35
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4338

Quant Time: Nov 15 22:27:21 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 : Wed Nov 13 00:41:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.549	152	2528	0.400	ng/ul	0.00
4) Naphthalene-d8	10.311	136	6318	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.184	164	5002	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.936	188	12349	0.400	ng/ul	0.00
17) Chrysene-d12	21.133	240	9837	0.400	ng/ul	0.00
23) Perylene-d12	23.296	264	8433	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	1006	0.418	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.911	152	4100	0.424	ng/ul	0.00
18) Fluoranthene-d10	18.967	212	12726	0.504	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.174	88	1077	0.409	ng/ul	97
5) Naphthalene	10.361	128	6648	0.420	ng/ul	98
7) 2-Methylnaphthalene	11.983	142	4953	0.420	ng/ul	100
8) 1-Methylnaphthalene	12.208	142	5074	0.428	ng/ul	99
10) Acenaphthylene	13.902	152	8103	0.397	ng/ul	99
11) Acenaphthene	14.249	153	6373	0.411	ng/ul	99
12) Fluorene	15.244	166	7936	0.403	ng/ul	98
14) Pentachlorophenol	16.594	266	991	0.347	ng/ul	98
15) Phenanthrene	16.979	178	13625	0.413	ng/ul	99
16) Anthracene	17.067	178	12468	0.402	ng/ul	98
19) Fluoranthene	19.000	202	16244	0.494	ng/ul#	96
20) Pyrene	19.362	202	16608	0.487	ng/ul#	95
21) Benzo(a)anthracene	21.116	228	14169	0.398	ng/ul	100
22) Chrysene	21.168	228	14421	0.397	ng/ul	100
24) Benzo(b)fluoranthene	22.653	252	12272	0.414	ng/ul	97
25) Benzo(k)fluoranthene	22.694	252	12911	0.416	ng/ul#	97
26) Benzo(a)pyrene	23.202	252	10928	0.413	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	25.443	276	13054	0.403	ng/ul#	92
28) Dibenzo(a,h)anthracene	25.459	278	10181	0.403	ng/ul#	97
29) Benzo(g,h,i)perylene	26.093	276	10183	0.411	ng/ul#	97

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

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