

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022822\
 Data File : BN018769.D
 Acq On : 28 Feb 2022 21:03
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4328

Quant Time: Mar 01 00:43:20 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN021822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 18 13:39:55 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.917	152	5532	0.400	ng/ul	0.00
4) Naphthalene-d8	10.721	136	13284	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.547	164	6179	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.283	188	11537	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.465	240	9608	0.400	ng/ul	# 0.00
23) Perylene-d12	23.864	264	8869	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	2727	0.434	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.310	152	7279	0.401	ng/ul	0.00
18) Fluoranthene-d10	19.312	212	11957	0.417	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.322	88	2961	0.460	ng/ul#	80
5) Naphthalene	10.770	128	15767	0.419	ng/ul	99
7) 2-Methylnaphthalene	12.381	142	9898	0.408	ng/ul	100
8) 1-Methylnaphthalene	12.596	142	9800	0.405	ng/ul	99
10) Acenaphthylene	14.265	152	10832	0.389	ng/ul	99
11) Acenaphthene	14.607	153	9443	0.423	ng/ul	99
12) Fluorene	15.592	166	10015	0.418	ng/ul	98
14) Pentachlorophenol	16.941	266	826	0.351	ng/ul	99
15) Phenanthrene	17.326	178	15958	0.421	ng/ul	99
16) Anthracene	17.419	178	13164	0.389	ng/ul	99
19) Fluoranthene	19.340	202	16954	0.420	ng/ul#	92
20) Pyrene	19.703	202	17453	0.419	ng/ul#	92
21) Benzo(a)anthracene	21.450	228	14365	0.386	ng/ul	99
22) Chrysene	21.502	228	16793	0.433	ng/ul	100
24) Benzo(b)fluoranthene	23.133	252	17527	0.444	ng/ul	93
25) Benzo(k)fluoranthene	23.180	252	17878	0.454	ng/ul#	94
26) Benzo(a)pyrene	23.756	252	14894	0.428	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	26.346	276	20066	0.453	ng/ul#	79
28) Dibenzo(a,h)anthracene	26.366	278	15421	0.443	ng/ul#	93
29) Benzo(g,h,i)perylene	27.111	276	16979	0.447	ng/ul#	91

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

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