

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072224\
 Data File : BN032897.D
 Acq On : 22 Jul 2024 17:23
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4274

Quant Time: Jul 22 17:55:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 20 01:52:15 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.532	152	3403	0.400	ng/ul	0.00
4) Naphthalene-d8	10.304	136	11219	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.182	164	5609	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.951	188	9704	0.400	ng/ul	0.00
17) Chrysene-d12	21.163	240	8908	0.400	ng/ul	# 0.00
23) Perylene-d12	23.358	264	11124	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.089	96	1602	0.392	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.909	152	6420	0.379	ng/ul	-0.02
18) Fluoranthene-d10	18.988	212	10233	0.352	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.122	88	1824	0.392	ng/ul#	79
5) Naphthalene	10.353	128	11492	0.381	ng/ul	99
7) 2-Methylnaphthalene	11.981	142	6954	0.370	ng/ul	99
8) 1-Methylnaphthalene	12.201	142	7074	0.364	ng/ul	99
10) Acenaphthylene	13.905	152	9729	0.429	ng/ul	99
11) Acenaphthene	14.242	153	7017	0.374	ng/ul	99
12) Fluorene	15.251	166	7364	0.344	ng/ul	99
14) Pentachlorophenol	16.635	266	912	0.536	ng/ul	97
15) Phenanthrene	16.993	178	9883	0.374	ng/ul	99
16) Anthracene	17.091	178	9160	0.424	ng/ul	98
19) Fluoranthene	19.021	202	12134	0.335	ng/ul	97
20) Pyrene	19.388	202	12910	0.344	ng/ul	96
21) Benzo(a)anthracene	21.149	228	12081	0.426	ng/ul	94
22) Chrysene	21.201	228	13538	0.330	ng/ul	94
24) Benzo(b)fluoranthene	22.697	252	13681	0.349	ng/ul	82
25) Benzo(k)fluoranthene	22.741	252	14137	0.296	ng/ul#	82
26) Benzo(a)pyrene	23.262	252	12034	0.354	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	25.564	276	17630	0.383	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.564	278	14056	0.394	ng/ul#	87
29) Benzo(g,h,i)perylene	26.245	276	13955	0.363	ng/ul#	89

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

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