

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072524\
 Data File : BN032966.D
 Acq On : 25 Jul 2024 10:22
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4280

Quant Time: Jul 26 01:07:04 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN072524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 25 15:43:00 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.506	152	2200	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.277	136	7356	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.165	164	5640	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.932	188	13217	0.400	ng/ul	-0.01
17) Chrysene-d12	21.153	240	14841	0.400	ng/ul	0.00
23) Perylene-d12	23.348	264	16530	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	977	0.351	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.878	152	4953	0.413	ng/ul	-0.03
18) Fluoranthene-d10	18.976	212	15586	0.371	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.110	88	1040	0.328	ng/ul#	45
5) Naphthalene	10.327	128	7448	0.394	ng/ul	96
7) 2-Methylnaphthalene	11.955	142	5302	0.414	ng/ul	95
8) 1-Methylnaphthalene	12.180	142	5483	0.396	ng/ul	99
10) Acenaphthylene	13.888	152	9457	0.410	ng/ul	99
11) Acenaphthene	14.230	153	6493	0.374	ng/ul	98
12) Fluorene	15.229	166	8026	0.397	ng/ul#	98
14) Pentachlorophenol	16.611	266	1724	0.641	ng/ul	98
15) Phenanthrene	16.974	178	13670	0.386	ng/ul	99
16) Anthracene	17.067	178	14001	0.439	ng/ul	99
19) Fluoranthene	19.008	202	17723	0.363	ng/ul	97
20) Pyrene	19.375	202	19155	0.366	ng/ul	99
21) Benzo(a)anthracene	21.138	228	20714	0.409	ng/ul	99
22) Chrysene	21.191	228	20751	0.351	ng/ul	99
24) Benzo(b)fluoranthene	22.690	252	21027	0.435	ng/ul	98
25) Benzo(k)fluoranthene	22.731	252	22543	0.390	ng/ul	97
26) Benzo(a)pyrene	23.254	252	17442	0.384	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	25.556	276	21748	0.315	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.552	278	17577	0.321	ng/ul#	93
29) Benzo(g,h,i)perylene	26.240	276	16672	0.304	ng/ul#	93

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

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