

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091424\
 Data File : BN033927.D
 Acq On : 14 Sep 2024 23:02
 Operator : JU/RC
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4305

Quant Time: Sep 15 02:00:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Sep 15 01:52:43 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.485	152	4383	0.400	ng/ul	0.00
4) Naphthalene-d8	10.244	136	12273	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.133	164	6197	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.885	188	12871	0.400	ng/ul	0.00
17) Chrysene-d12	21.100	240	11140	0.400	ng/ul	0.00
23) Perylene-d12	23.245	264	12473	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.101	96	1992	0.442	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.845	152	6827	0.388	ng/ul	0.00
18) Fluoranthene-d10	18.924	212	13058	0.393	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.135	88	2222	0.444	ng/ul#	86
5) Naphthalene	10.294	128	12806	0.390	ng/ul	100
7) 2-Methylnaphthalene	11.916	142	7991	0.384	ng/ul	100
8) 1-Methylnaphthalene	12.142	142	8335	0.390	ng/ul	99
10) Acenaphthylene	13.846	152	10758	0.382	ng/ul	99
11) Acenaphthene	14.193	153	8305	0.409	ng/ul	98
12) Fluorene	15.187	166	9260	0.408	ng/ul	100
14) Pentachlorophenol	16.543	266	1209	0.299	ng/ul	99
15) Phenanthrene	16.927	178	14574	0.384	ng/ul	99
16) Anthracene	17.016	178	12494	0.366	ng/ul	100
19) Fluoranthene	18.957	202	16582	0.377	ng/ul	99
20) Pyrene	19.319	202	17395	0.373	ng/ul	99
21) Benzo(a)anthracene	21.086	228	15505	0.354	ng/ul	99
22) Chrysene	21.138	228	17917	0.404	ng/ul	100
24) Benzo(b)fluoranthene	22.608	252	17590	0.363	ng/ul	98
25) Benzo(k)fluoranthene	22.652	252	19730	0.407	ng/ul	98
26) Benzo(a)pyrene	23.152	252	14770	0.380	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.361	276	22698	0.389	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.378	278	17509	0.386	ng/ul	97
29) Benzo(g,h,i)perylene	26.002	276	18220	0.399	ng/ul	99

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

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