

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091624\
 Data File : BN033999.D
 Acq On : 16 Sep 2024 22:26
 Operator : JU/RC
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4311

Quant Time: Sep 17 00:42:39 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	7.477	152	6170	0.400	ng/ul	0.00
4) Naphthalene-d8	10.233	136	11327	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.119	164	5510	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.877	188	10875	0.400	ng/ul	0.00
17) Chrysene-d12	21.095	240	8785	0.400	ng/ul	0.00
23) Perylene-d12	23.234	264	9784	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.097	96	2056	0.324	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.834	152	6481	0.399	ng/ul	-0.01
18) Fluoranthene-d10	18.915	212	11447	0.437	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.131	88	2251	0.320	ng/ul#	89
5) Naphthalene	10.283	128	12365	0.408	ng/ul	100
7) 2-Methylnaphthalene	11.911	142	7603	0.396	ng/ul	99
8) 1-Methylnaphthalene	12.131	142	7842	0.397	ng/ul	100
10) Acenaphthylene	13.837	152	9866	0.394	ng/ul	99
11) Acenaphthene	14.184	153	7698	0.426	ng/ul	99
12) Fluorene	15.178	166	8403	0.416	ng/ul	100
14) Pentachlorophenol	16.539	266	977	0.286	ng/ul	100
15) Phenanthrene	16.919	178	12756	0.398	ng/ul	100
16) Anthracene	17.012	178	10907	0.378	ng/ul	100
19) Fluoranthene	18.948	202	14474	0.418	ng/ul	99
20) Pyrene	19.310	202	15197	0.413	ng/ul	100
21) Benzo(a)anthracene	21.077	228	13242	0.383	ng/ul	99
22) Chrysene	21.127	228	15047	0.431	ng/ul	99
24) Benzo(b)fluoranthene	22.599	252	14678	0.386	ng/ul	98
25) Benzo(k)fluoranthene	22.640	252	16310	0.429	ng/ul	99
26) Benzo(a)pyrene	23.140	252	12343	0.405	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.348	276	18953	0.414	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.365	278	14652	0.412	ng/ul	98
29) Benzo(g,h,i)perylene	25.992	276	15268	0.427	ng/ul	99

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

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