

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121624\
 Data File : BN035615.D
 Acq On : 16 Dec 2024 14:02
 Operator : RC/JU
 Sample : SSTD0.445
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4232

Quant Time: Dec 16 14:42:54 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN121624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 16 14:42:43 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.284	152	1866	0.400	ng/ul	0.00
4) Naphthalene-d8	10.031	136	5515	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.949	164	3919	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.717	188	9201	0.400	ng/ul	0.00
17) Chrysene-d12	20.964	240	8386	0.400	ng/ul	0.00
23) Perylene-d12	23.054	264	9372	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.955	96	706	0.418	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.637	152	3176	0.398	ng/ul	0.00
18) Fluoranthene-d10	18.767	212	9775	0.422	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.989	88	765	0.416	ng/ul	100
5) Naphthalene	10.080	128	5971	0.428	ng/ul	100
7) 2-Methylnaphthalene	11.714	142	3691	0.391	ng/ul	100
8) 1-Methylnaphthalene	11.939	142	3824	0.398	ng/ul	100
10) Acenaphthylene	13.662	152	6248	0.404	ng/ul	100
11) Acenaphthene	14.009	153	4725	0.400	ng/ul	100
12) Fluorene	15.013	166	5532	0.400	ng/ul	100
14) Pentachlorophenol	16.379	266	889	0.406	ng/ul	100
15) Phenanthrene	16.759	178	9341	0.389	ng/ul	100
16) Anthracene	16.848	178	8721	0.391	ng/ul	100
19) Fluoranthene	18.800	202	11913	0.411	ng/ul	100
20) Pyrene	19.167	202	13109	0.430	ng/ul	100
21) Benzo(a)anthracene	20.950	228	11262	0.393	ng/ul	100
22) Chrysene	20.999	228	11868	0.408	ng/ul	100
24) Benzo(b)fluoranthene	22.442	252	11377	0.376	ng/ul	100
25) Benzo(k)fluoranthene	22.480	252	12269	0.391	ng/ul	100
26) Benzo(a)pyrene	22.963	252	10798	0.387	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.081	276	13406	0.363	ng/ul	100
28) Dibenzo(a,h)anthracene	25.101	278	10297	0.357	ng/ul	100
29) Benzo(g,h,i)perylene	25.698	276	10643	0.365	ng/ul	100

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

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