

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112624\
 Data File : BN035330.D
 Acq On : 27 Nov 2024 00:18
 Operator : RC/JU
 Sample : P4996-03MSD
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHF70MSD

Quant Time: Nov 27 01:00:28 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN112624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 06:09:15 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.310	152	3473	0.400	ng/ul	0.00
4) Naphthalene-d8	10.059	136	9177	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.967	164	5537	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.734	188	10850	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.976	240	6359	0.400	ng/ul	# 0.00
23) Perylene-d12	23.068	264	6851	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.976	96	688	0.217	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.659	152	3434	0.255	ng/ul	0.00
18) Fluoranthene-d10	18.786	212	8124	0.463	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.006	88	2086	0.605	ng/ul	93
5) Naphthalene	10.108	128	6234	0.270	ng/ul	100
7) 2-Methylnaphthalene	11.736	142	4143	0.258	ng/ul	99
8) 1-Methylnaphthalene	11.961	142	4300	0.266	ng/ul	99
10) Acenaphthylene	13.681	152	6756	0.308	ng/ul	99
11) Acenaphthene	14.032	153	5089	0.303	ng/ul	98
12) Fluorene	15.036	166	5975	0.303	ng/ul	99
14) Pentachlorophenol	16.392	266	1779	0.704	ng/ul	98
15) Phenanthrene	16.776	178	10027	0.350	ng/ul	100
16) Anthracene	16.869	178	8516	0.320	ng/ul	99
19) Fluoranthene	18.819	202	10105	0.456	ng/ul	97
20) Pyrene	19.181	202	10102	0.438	ng/ul#	91
21) Benzo(a)anthracene	20.958	228	7869	0.357	ng/ul	99
22) Chrysene	21.014	228	8539	0.385	ng/ul	99
24) Benzo(b)fluoranthene	22.454	252	8146	0.360	ng/ul	96
25) Benzo(k)fluoranthene	22.498	252	9146	0.392	ng/ul#	96
26) Benzo(a)pyrene	22.978	252	7320	0.354	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	25.108	276	10222	0.366	ng/ul#	78
28) Dibenzo(a,h)anthracene	25.128	278	7719	0.352	ng/ul#	95
29) Benzo(g,h,i)perylene	25.725	276	8157	0.370	ng/ul#	94

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

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