

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121024\
 Data File : BN035565.D
 Acq On : 11 Dec 2024 14:16
 Operator : RC/JU
 Sample : P5152-03MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHK72MSD

Quant Time: Dec 11 14:49:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 10 16:10:57 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.293	152	2424	0.400	ng/ul	0.00
4) Naphthalene-d8	10.036	136	6849	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.949	164	4651	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.717	188	10145	0.400	ng/ul	0.00
17) Chrysene-d12	20.964	240	8263	0.400	ng/ul	0.00
23) Perylene-d12	23.053	264	9170	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.963	96	613	0.277	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.637	152	3547	0.353	ng/ul	0.00
18) Fluoranthene-d10	18.772	212	9957	0.437	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.993	88	1781	0.741	ng/ul#	86
5) Naphthalene	10.086	128	6804	0.394	ng/ul	100
7) 2-Methylnaphthalene	11.714	142	3966	0.331	ng/ul	97
8) 1-Methylnaphthalene	11.939	142	4058	0.336	ng/ul	99
10) Acenaphthylene	13.662	152	6086	0.330	ng/ul	99
11) Acenaphthene	14.013	153	4741	0.336	ng/ul	99
12) Fluorene	15.013	166	5567	0.336	ng/ul	99
14) Pentachlorophenol	16.379	266	959	0.406	ng/ul	98
15) Phenanthrene	16.759	178	9729	0.363	ng/ul	100
16) Anthracene	16.852	178	8618	0.346	ng/ul	99
19) Fluoranthene	18.800	202	11785	0.410	ng/ul	98
20) Pyrene	19.167	202	12619	0.421	ng/ul#	96
21) Benzo(a)anthracene	20.947	228	11614	0.405	ng/ul	99
22) Chrysene	20.999	228	11948	0.414	ng/ul	100
24) Benzo(b)fluoranthene	22.439	252	11346	0.374	ng/ul	97
25) Benzo(k)fluoranthene	22.483	252	12038	0.385	ng/ul	97
26) Benzo(a)pyrene	22.963	252	10858	0.392	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	25.084	276	14030	0.375	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.104	278	10049	0.343	ng/ul#	94
29) Benzo(g,h,i)perylene	25.701	276	10156	0.344	ng/ul#	95

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

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