

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN103124\
 Data File : BN034772.D
 Acq On : 31 Oct 2024 18:53
 Operator : RC/JU
 Sample : P4435-03MSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHK68MSD

Quant Time: Nov 01 05:02:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 31 11:13:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.604	152	6919	0.400	ng/ul	0.00
4) Naphthalene-d8	10.367	136	21471	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.236	164	11736	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.983	188	24021	0.400	ng/ul	0.00
17) Chrysene-d12	21.175	240	17440	0.400	ng/ul	0.00
23) Perylene-d12	23.355	264	15642	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.174	96	2615	0.365	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.967	152	6355	0.236	ng/ul	0.00
18) Fluoranthene-d10	19.014	212	15054	0.265	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.204	88	7446	0.947	ng/ul#	88
5) Naphthalene	10.416	128	12848	0.226	ng/ul	98
7) 2-Methylnaphthalene	12.038	142	8489	0.230	ng/ul	100
8) 1-Methylnaphthalene	12.258	142	10100	0.268	ng/ul	100
10) Acenaphthylene	13.954	152	14599	0.273	ng/ul	99
11) Acenaphthene	14.300	153	9994	0.245	ng/ul	98
12) Fluorene	15.291	166	11566	0.254	ng/ul	98
14) Pentachlorophenol	16.633	266	764	0.161	ng/ul	99
15) Phenanthrene	17.021	178	30862	0.425	ng/ul	99
16) Anthracene	17.114	178	17846	0.265	ng/ul	99
19) Fluoranthene	19.042	202	41406	0.489	ng/ul	97
20) Pyrene	19.404	202	43303	0.492	ng/ul	97
21) Benzo(a)anthracene	21.157	228	26350	0.378	ng/ul	98
22) Chrysene	21.210	228	27806	0.398	ng/ul	99
24) Benzo(b)fluoranthene	22.706	252	26095	0.410	ng/ul	98
25) Benzo(k)fluoranthene	22.750	252	21073	0.328	ng/ul	98
26) Benzo(a)pyrene	23.258	252	20802	0.380	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.530	276	19222	0.329	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.547	278	13092	0.297	ng/ul	99
29) Benzo(g,h,i)perylene	26.188	276	15864	0.346	ng/ul	100

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

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