

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110923\
 Data File : BN028547.D
 Acq On : 10 Nov 2023 08:57
 Operator : MA/JU
 Sample : 05044-03MSD
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHEE3MSD

Quant Time: Nov 10 23:25:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 09 16:30:32 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.703	152	6484	0.400	ng/ul	0.00
4) Naphthalene-d8	10.491	136	21719	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.349	164	13658	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.099	188	30461	0.400	ng/ul	0.00
17) Chrysene-d12	21.312	240	24769	0.400	ng/ul	0.00
23) Perylene-d12	23.618	264	23204	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.151	96	2070	0.289	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.091	152	11491	0.351	ng/ul	0.00
18) Fluoranthene-d10	19.137	212	38370	0.481	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.185	88	9161	1.157	ng/ul#	84
5) Naphthalene	10.540	128	22806	0.375	ng/ul	97
7) 2-Methylnaphthalene	12.163	142	15104	0.377	ng/ul	97
8) 1-Methylnaphthalene	12.383	142	15989	0.393	ng/ul	98
10) Acenaphthylene	14.067	152	26266	0.385	ng/ul	97
11) Acenaphthene	14.414	153	19591	0.391	ng/ul	98
12) Fluorene	15.404	166	24038	0.412	ng/ul	99
14) Pentachlorophenol	16.753	266	168	0.023	ng/ul#	85
15) Phenanthrene	17.141	178	46593	0.498	ng/ul	99
16) Anthracene	17.234	178	41573	0.484	ng/ul	98
19) Fluoranthene	19.169	202	54085	0.513	ng/ul	98
20) Pyrene	19.532	202	63383	0.581	ng/ul	98
21) Benzo(a)anthracene	21.298	228	49186	0.502	ng/ul	99
22) Chrysene	21.347	228	49156	0.513	ng/ul	99
24) Benzo(b)fluoranthene	22.917	252	43641	0.501	ng/ul	100
25) Benzo(k)fluoranthene	22.966	252	44490	0.488	ng/ul	98
26) Benzo(a)pyrene	23.513	252	39054	0.475	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.980	276	42115	0.462	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.003	278	34324	0.471	ng/ul	98
29) Benzo(g,h,i)perylene	26.701	276	34356	0.459	ng/ul	99

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

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