

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110923\
 Data File : BN028546.D
 Acq On : 10 Nov 2023 08:21
 Operator : MA/JU
 Sample : 05044-02MS
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHEE3MS

Quant Time: Nov 10 23:25:44 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	7426	0.400	ng/ul	0.00
4) Naphthalene-d8	10.489	136	24656	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.352	164	15550	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.102	188	34714	0.400	ng/ul	0.00
17) Chrysene-d12	21.315	240	27979	0.400	ng/ul	0.00
23) Perylene-d12	23.616	264	25789	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.151	96	2177	0.265	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.090	152	11874	0.320	ng/ul	0.00
18) Fluoranthene-d10	19.140	212	39357	0.437	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.184	88	9441	1.041	ng/ul#	84
5) Naphthalene	10.539	128	23523	0.340	ng/ul	98
7) 2-Methylnaphthalene	12.161	142	15592	0.343	ng/ul	97
8) 1-Methylnaphthalene	12.381	142	16483	0.357	ng/ul	99
10) Acenaphthylene	14.070	152	26846	0.346	ng/ul	97
11) Acenaphthene	14.413	153	20475	0.359	ng/ul	99
12) Fluorene	15.403	166	24821	0.374	ng/ul	99
14) Pentachlorophenol	16.751	266	230	0.028	ng/ul	97
15) Phenanthrene	17.140	178	48127	0.451	ng/ul	98
16) Anthracene	17.233	178	43171	0.441	ng/ul	98
19) Fluoranthene	19.168	202	55415	0.466	ng/ul	99
20) Pyrene	19.535	202	65219	0.529	ng/ul	99
21) Benzo(a)anthracene	21.298	228	50010	0.452	ng/ul	98
22) Chrysene	21.350	228	49896	0.461	ng/ul	99
24) Benzo(b)fluoranthene	22.920	252	43706	0.452	ng/ul	100
25) Benzo(k)fluoranthene	22.967	252	44940	0.443	ng/ul	100
26) Benzo(a)pyrene	23.513	252	38934	0.426	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.980	276	41199	0.406	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.000	278	33520	0.414	ng/ul	97
29) Benzo(g,h,i)perylene	26.701	276	33188	0.399	ng/ul	99

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

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