

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112224\
 Data File : BN035261.D
 Acq On : 23 Nov 2024 03:23
 Operator : RC/JU
 Sample : P4872-06MS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E29X5MS

Quant Time: Nov 23 04:05:56 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 : Fri Nov 22 12:45:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.318	152	4409	0.400	ng/ul	0.00
4) Naphthalene-d8	10.069	136	11409	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.977	164	7219	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.747	188	15229	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.988	240	10345	0.400	ng/ul	# 0.00
23) Perylene-d12	23.089	264	9174	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.976	96	1191	0.296	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.670	152	3911	0.234	ng/ul	0.00
18) Fluoranthene-d10	18.795	212	9764	0.342	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.005	88	3562	0.814	ng/ul#	81
5) Naphthalene	10.113	128	7015	0.244	ng/ul	99
7) 2-Methylnaphthalene	11.747	142	4838	0.243	ng/ul	98
8) 1-Methylnaphthalene	11.972	142	4928	0.245	ng/ul	98
10) Acenaphthylene	13.690	152	8461	0.295	ng/ul	99
11) Acenaphthene	14.041	153	5759	0.263	ng/ul	97
12) Fluorene	15.045	166	6687	0.260	ng/ul	99
14) Pentachlorophenol	16.405	266	1505	0.424	ng/ul	98
15) Phenanthrene	16.785	178	14488	0.360	ng/ul	99
16) Anthracene	16.878	178	10850	0.290	ng/ul	100
19) Fluoranthene	18.828	202	24415	0.678	ng/ul#	94
20) Pyrene	19.195	202	22464	0.599	ng/ul#	92
21) Benzo(a)anthracene	20.973	228	14449	0.403	ng/ul	98
22) Chrysene	21.026	228	14413	0.399	ng/ul	99
24) Benzo(b)fluoranthene	22.475	252	13281	0.438	ng/ul	96
25) Benzo(k)fluoranthene	22.513	252	11048	0.353	ng/ul#	95
26) Benzo(a)pyrene	22.998	252	10712	0.387	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.138	276	12769	0.341	ng/ul#	84
28) Dibenzo(a,h)anthracene	25.155	278	8692	0.296	ng/ul#	92
29) Benzo(g,h,i)perylene	25.755	276	10325	0.350	ng/ul#	92

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

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