

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092823\
 Data File : BN027924.D
 Acq On : 28 Sep 2023 14:03
 Operator : MA/JU
 Sample : 04417-39
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EZYE6

Quant Time: Sep 28 16:27:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN091523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 27 03:52:19 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	3463	0.400	ng/ul	0.00
4) Naphthalene-d8	10.779	136	12971	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.601	164	9162	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.354	188	18631	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.551	240	15545	0.400	ng/ul	# 0.01
23) Perylene-d12	24.032	264	13454	0.400	ng/ul	# 0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.325	96	23234	5.326	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.363	152	10688	0.544	ng/ul	0.00
18) Fluoranthene-d10	19.385	212	23658	0.498	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.372	88	102	0.022	ng/ul#	1
5) Naphthalene	10.828	128	60081	1.621	ng/ul	100
7) 2-Methylnaphthalene	12.434	142	36467	1.550	ng/ul	100
8) 1-Methylnaphthalene	12.654	142	25600	1.055	ng/ul	99
10) Acenaphthylene	14.333	152	8164	0.204	ng/ul#	57
11) Acenaphthene	14.670	153	2983	0.086	ng/ul#	69
12) Fluorene	15.651	166	2416	0.059	ng/ul#	24
14) Pentachlorophenol	16.987	266	267	0.061	ng/ul#	50
15) Phenanthrene	17.401	178	352104	6.014	ng/ul	97
16) Anthracene	17.489	178	24131	0.492	ng/ul#	69
19) Fluoranthene	19.418	202	670116	10.660	ng/ul	100
20) Pyrene	19.785	202	580471	8.964	ng/ul	100
21) Benzo(a)anthracene	21.533	228	304576	5.064	ng/ul	96
22) Chrysene	21.586	228	462270	7.438	ng/ul	96
24) Benzo(b)fluoranthene	23.266	252	809004	13.490	ng/ul	86
26) Benzo(a)pyrene	23.918	252	273773	5.676	ng/ul#	81
27) Indeno(1,2,3-cd)pyrene	26.633	276	295107	4.775	ng/ul#	86
28) Dibenzo(a,h)anthracene	26.643	278	85790	1.724	ng/ul#	60
29) Benzo(g,h,i)perylene	27.448	276	275984	5.258	ng/ul	99

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

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