

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112224\
 Data File : BN035256.D
 Acq On : 23 Nov 2024 00:22
 Operator : RC/JU
 Sample : PB165099BS
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS099

Quant Time: Nov 23 00:51:37 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	3718	0.400	ng/ul	0.00
4) Naphthalene-d8	10.069	136	9412	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.977	164	5930	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.747	188	13279	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.988	240	9950	0.400	ng/ul	# 0.00
23) Perylene-d12	23.092	264	8640	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.976	96	1723	0.507	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.670	152	4211	0.305	ng/ul	0.00
18) Fluoranthene-d10	18.795	212	10277	0.374	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.005	88	4894	1.327	ng/ul#	89
5) Naphthalene	10.113	128	7549	0.318	ng/ul	99
7) 2-Methylnaphthalene	11.747	142	5106	0.310	ng/ul	97
8) 1-Methylnaphthalene	11.972	142	5217	0.314	ng/ul	99
10) Acenaphthylene	13.690	152	8117	0.345	ng/ul	100
11) Acenaphthene	14.041	153	6028	0.335	ng/ul	97
12) Fluorene	15.045	166	6877	0.326	ng/ul	99
14) Pentachlorophenol	16.405	266	2052	0.663	ng/ul	99
15) Phenanthrene	16.789	178	11522	0.328	ng/ul	99
16) Anthracene	16.878	178	10250	0.315	ng/ul	100
19) Fluoranthene	18.828	202	13291	0.384	ng/ul#	94
20) Pyrene	19.195	202	13612	0.378	ng/ul#	91
21) Benzo(a)anthracene	20.973	228	11117	0.322	ng/ul	99
22) Chrysene	21.023	228	11742	0.338	ng/ul	100
24) Benzo(b)fluoranthene	22.475	252	9660	0.338	ng/ul	95
25) Benzo(k)fluoranthene	22.519	252	10576	0.359	ng/ul#	95
26) Benzo(a)pyrene	23.001	252	8563	0.328	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.141	276	10525	0.299	ng/ul#	83
28) Dibenzo(a,h)anthracene	25.158	278	7921	0.287	ng/ul#	92
29) Benzo(g,h,i)perylene	25.762	276	8291	0.298	ng/ul#	93

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

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