

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092623\
 Data File : BN027871.D
 Acq On : 26 Sep 2023 20:57
 Operator : MA/JU
 Sample : PB155614BS
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS614

Quant Time: Sep 27 03:01:42 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 : Fri Sep 22 03:58:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.966	152	2224	0.400	ng/ul	0.00
4) Naphthalene-d8	10.784	136	7803	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.606	164	5356	0.400	ng/ul	0.01
13) Phenanthrene-d10	17.358	188	13008	0.400	ng/ul	0.02
17) Chrysene-d12	21.542	240	13388	0.400	ng/ul	0.02
23) Perylene-d12	24.015	264	15286	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	2209	0.789	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.368	152	5411	0.458	ng/ul	0.01
18) Fluoranthene-d10	19.380	212	15905	0.389	ng/ul	0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.368	88	6665	2.286	ng/ul#	86
5) Naphthalene	10.834	128	8531	0.383	ng/ul	98
7) 2-Methylnaphthalene	12.440	142	5797	0.410	ng/ul	100
8) 1-Methylnaphthalene	12.660	142	5719	0.392	ng/ul	99
10) Acenaphthylene	14.333	152	9089	0.388	ng/ul	100
11) Acenaphthene	14.670	153	7221	0.357	ng/ul	98
12) Fluorene	15.656	166	8625	0.363	ng/ul	100
14) Pentachlorophenol	16.987	266	2702	0.888	ng/ul	98
15) Phenanthrene	17.401	178	14831	0.363	ng/ul	99
16) Anthracene	17.494	178	12887	0.377	ng/ul	99
19) Fluoranthene	19.413	202	18033	0.333	ng/ul	99
20) Pyrene	19.780	202	18954	0.340	ng/ul	99
21) Benzo(a)anthracene	21.524	228	19832	0.383	ng/ul	100
22) Chrysene	21.580	228	19859	0.371	ng/ul	99
24) Benzo(b)fluoranthene	23.249	252	21889	0.321	ng/ul	96
25) Benzo(k)fluoranthene	23.296	252	22466	0.340	ng/ul	96
26) Benzo(a)pyrene	23.901	252	21049	0.384	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.613	276	27234	0.388	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.636	278	22281	0.394	ng/ul	98
29) Benzo(g,h,i)perylene	27.425	276	22729	0.381	ng/ul	98

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

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