

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082322\
 Data File : BN021328.D
 Acq On : 23 Aug 2022 14:38
 Operator : CG/JU
 Sample : PB147164BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS164

Quant Time: Aug 24 00:58:42 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 22 18:06:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.097	152	2742	0.400	ng/ul	0.00
4) Naphthalene-d8	10.927	136	8598	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.749	164	4668	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.498	188	9850	0.400	ng/ul	0.00
17) Chrysene-d12	21.691	240	8216	0.400	ng/ul	# 0.00
23) Perylene-d12	24.225	264	6831	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.384	96	2609	0.738	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.511	152	5354	0.412	ng/ul	0.00
18) Fluoranthene-d10	19.524	212	10631	0.420	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.418	88	6837	1.845	ng/ul#	87
5) Naphthalene	10.977	128	9560	0.351	ng/ul	98
7) 2-Methylnaphthalene	12.582	142	6120	0.359	ng/ul	98
8) 1-Methylnaphthalene	12.802	142	6372	0.399	ng/ul	99
10) Acenaphthylene	14.471	152	8203	0.340	ng/ul	98
11) Acenaphthene	14.809	153	6448	0.350	ng/ul	97
12) Fluorene	15.794	166	7184	0.349	ng/ul	99
14) Pentachlorophenol	17.139	266	846	0.778	ng/ul	99
15) Phenanthrene	17.540	178	12608	0.345	ng/ul	100
16) Anthracene	17.633	178	9994	0.342	ng/ul	99
19) Fluoranthene	19.557	202	13851	0.356	ng/ul	99
20) Pyrene	19.919	202	13562	0.348	ng/ul	98
21) Benzo(a)anthracene	21.673	228	10786	0.379	ng/ul	99
22) Chrysene	21.729	228	13595	0.371	ng/ul	99
24) Benzo(b)fluoranthene	23.445	252	12564	0.428	ng/ul	98
25) Benzo(k)fluoranthene	23.497	252	12638	0.400	ng/ul	96
26) Benzo(a)pyrene	24.111	252	10743	0.389	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.901	276	13211	0.407	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.928	278	11115	0.416	ng/ul	95
29) Benzo(g,h,i)perylene	27.726	276	12714	0.394	ng/ul	99

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

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