

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081522\
 Data File : BN021234.D
 Acq On : 15 Aug 2022 13:43
 Operator : CG/JU
 Sample : PB146922BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS922

Quant Time: Aug 15 14:28:35 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 13 04:02:44 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.797	152	2720	0.400	ng/u1	0.00
4) Naphthalene-d8	10.580	136	9896	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.433	164	5652	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.188	188	9994	0.400	ng/u1	0.00
17) Chrysene-d12	21.395	240	6412	0.400	ng/u1	0.00
23) Perylene-d12	23.727	264	5130	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.194	96	2395	0.688	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.186	152	4967	0.343	ng/u1	0.00
18) Fluoranthene-d10	19.226	212	9430	0.476	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.228	88	6132	1.711	ng/u1	86
5) Naphthalene	10.635	128	9413	0.310	ng/u1	99
7) 2-Methylnaphthalene	12.263	142	5582	0.303	ng/u1	98
8) 1-Methylnaphthalene	12.477	142	6150	0.333	ng/u1	100
10) Acenaphthylene	14.156	152	8020	0.303	ng/u1	99
11) Acenaphthene	14.498	153	6591	0.302	ng/u1	99
12) Fluorene	15.493	166	7011	0.293	ng/u1	100
14) Pentachlorophenol	16.859	266	574	0.372	ng/u1	97
15) Phenanthrene	17.231	178	10897	0.305	ng/u1	100
16) Anthracene	17.328	178	8366	0.302	ng/u1	100
19) Fluoranthene	19.259	202	11553	0.407	ng/u1	100
20) Pyrene	19.621	202	11994	0.409	ng/u1	97
21) Benzo(a)anthracene	21.383	228	6302	0.301	ng/u1	98
22) Chrysene	21.433	228	8602	0.312	ng/u1	99
24) Benzo(b)fluoranthene	23.020	252	7666	0.323	ng/u1	88
25) Benzo(k)fluoranthene	23.067	252	8056	0.321	ng/u1	98
26) Benzo(a)pyrene	23.625	252	7425	0.343	ng/u1#	69
27) Indeno(1,2,3-cd)pyrene	26.139	276	7730	0.299	ng/u1	94
28) Dibenzo(a,h)anthracene	26.155	278	5724	0.280	ng/u1	96
29) Benzo(g,h,i)perylene	26.866	276	7092	0.307	ng/u1	98

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

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