

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121322\
 Data File : BN023195.D
 Acq On : 14 Dec 2022 10:46
 Operator : CG/JU
 Sample : PB149540BS
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS540

Quant Time: Dec 14 23:19:22 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN112622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 13 23:52:47 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.015	152	4608	0.400	ng/u1	0.00
4) Naphthalene-d8	10.832	136	14450	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.659	164	8393	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.403	188	19705	0.400	ng/u1	0.00
17) Chrysene-d12	21.587	240	19299	0.400	ng/u1	0.00
23) Perylene-d12	24.039	264	16147	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	3094	0.560	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.416	152	7097	0.322	ng/u1	0.00
18) Fluoranthene-d10	19.430	212	17714	0.248	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.391	88	7902	1.370	ng/u1	93
5) Naphthalene	10.881	128	13104	0.299	ng/u1	99
7) 2-Methylnaphthalene	12.493	142	8329	0.304	ng/u1	100
8) 1-Methylnaphthalene	12.707	142	8951	0.321	ng/u1	100
10) Acenaphthylene	14.382	152	11081	0.268	ng/u1	100
11) Acenaphthene	14.719	153	9709	0.296	ng/u1	98
12) Fluorene	15.705	166	11528	0.311	ng/u1	99
14) Pentachlorophenol	17.052	266	3034	0.736	ng/u1#	100
15) Phenanthrene	17.445	178	20384	0.301	ng/u1	100
16) Anthracene	17.538	178	16503	0.284	ng/u1	99
19) Fluoranthene	19.462	202	24163	0.242	ng/u1	99
20) Pyrene	19.820	202	24470	0.250	ng/u1	96
21) Benzo(a)anthracene	21.569	228	24045	0.326	ng/u1	100
22) Chrysene	21.625	228	26464	0.327	ng/u1	99
24) Benzo(b)fluoranthene	23.288	252	27567	0.337	ng/u1	96
25) Benzo(k)fluoranthene	23.338	252	25893	0.299	ng/u1	97
26) Benzo(a)pyrene	23.928	252	22251	0.357	ng/u1#	93
27) Indeno(1,2,3-cd)pyrene	26.604	276	29297	0.396	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.627	278	23028	0.391	ng/u1	99
29) Benzo(g,h,i)perylene	27.395	276	23955	0.371	ng/u1	99

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

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