

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121322\
 Data File : BN023181.D
 Acq On : 14 Dec 2022 01:39
 Operator : CG/JU
 Sample : PB149489BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS489

Quant Time: Dec 14 03:28:18 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	6391	0.400	ng/u1	0.00
4) Naphthalene-d8	10.832	136	19645	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.659	164	11372	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.403	188	25970	0.400	ng/u1	0.00
17) Chrysene-d12	21.588	240	24493	0.400	ng/u1	0.00
23) Perylene-d12	24.044	264	19527	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	4218	0.550	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.421	152	9192	0.307	ng/u1	0.00
18) Fluoranthene-d10	19.430	212	21970	0.243	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.395	88	11283	1.411	ng/u1	94
5) Naphthalene	10.881	128	17853	0.299	ng/u1	100
7) 2-Methylnaphthalene	12.493	142	11344	0.304	ng/u1	100
8) 1-Methylnaphthalene	12.712	142	12232	0.323	ng/u1	100
10) Acenaphthylene	14.382	152	14774	0.263	ng/u1	99
11) Acenaphthene	14.724	153	13094	0.295	ng/u1	99
12) Fluorene	15.705	166	15175	0.302	ng/u1	99
14) Pentachlorophenol	17.052	266	3572	0.658	ng/u1#	100
15) Phenanthrene	17.445	178	26920	0.302	ng/u1	100
16) Anthracene	17.538	178	21742	0.284	ng/u1	100
19) Fluoranthene	19.462	202	31611	0.250	ng/u1	99
20) Pyrene	19.825	202	31806	0.257	ng/u1	99
21) Benzo(a)anthracene	21.571	228	30377	0.324	ng/u1	100
22) Chrysene	21.623	228	33689	0.328	ng/u1	99
24) Benzo(b)fluoranthene	23.290	252	33329	0.337	ng/u1	98
25) Benzo(k)fluoranthene	23.336	252	32603	0.312	ng/u1	99
26) Benzo(a)pyrene	23.930	252	27137	0.360	ng/u1#	96
27) Indeno(1,2,3-cd)pyrene	26.609	276	34644	0.387	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.629	278	27298	0.384	ng/u1	98
29) Benzo(g,h,i)perylene	27.397	276	28176	0.361	ng/u1	99

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

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