

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091424\
 Data File : BN033956.D
 Acq On : 15 Sep 2024 17:39
 Operator : JU/RC
 Sample : PB163194BS
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS194

Quant Time: Sep 16 00:14:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Sep 15 01:52:43 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.481	152	4496	0.400	ng/ul	0.00
4) Naphthalene-d8	10.239	136	12490	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.124	164	6344	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.881	188	13318	0.400	ng/ul	0.00
17) Chrysene-d12	21.095	240	11249	0.400	ng/ul	0.00
23) Perylene-d12	23.237	264	12226	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.097	96	3863	0.836	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.839	152	7120	0.397	ng/ul	0.00
18) Fluoranthene-d10	18.920	212	15113	0.451	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.131	88	10111	1.971	ng/ul#	79
5) Naphthalene	10.288	128	12122	0.363	ng/ul	100
7) 2-Methylnaphthalene	11.916	142	7537	0.356	ng/ul	99
8) 1-Methylnaphthalene	12.136	142	9066	0.417	ng/ul	100
10) Acenaphthylene	13.841	152	9789	0.340	ng/ul	99
11) Acenaphthene	14.188	153	8065	0.388	ng/ul	98
12) Fluorene	15.183	166	8646	0.372	ng/ul	99
14) Pentachlorophenol	16.543	266	739	0.176	ng/ul	97
15) Phenanthrene	16.923	178	13942	0.355	ng/ul	100
16) Anthracene	17.016	178	12086	0.342	ng/ul	100
19) Fluoranthene	18.953	202	16593	0.374	ng/ul	100
20) Pyrene	19.315	202	17581	0.373	ng/ul	100
21) Benzo(a)anthracene	21.080	228	15317	0.346	ng/ul	100
22) Chrysene	21.130	228	17761	0.397	ng/ul	99
24) Benzo(b)fluoranthene	22.602	252	16795	0.354	ng/ul	98
25) Benzo(k)fluoranthene	22.646	252	18806	0.396	ng/ul	98
26) Benzo(a)pyrene	23.143	252	15599	0.409	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.355	276	21142	0.369	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.368	278	16283	0.366	ng/ul	98
29) Benzo(g,h,i)perylene	25.995	276	17225	0.385	ng/ul	98

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

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