

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100523\
 Data File : BN028109.D
 Acq On : 07 Oct 2023 06:53
 Operator : MA/JU
 Sample : PB156049BS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS049

Quant Time: Oct 08 00:50:17 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN100323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 06 05:15:40 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.933	152	4547	0.400	ng/ul	0.00
4) Naphthalene-d8	10.751	136	14149	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.578	164	7051	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.333	188	13332	0.400	ng/ul	0.00
17) Chrysene-d12	21.522	240	11978	0.400	ng/ul	0.00
23) Perylene-d12	23.986	264	13345	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.326	96	4639	0.812	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.335	152	8897	0.390	ng/ul	0.00
18) Fluoranthene-d10	19.362	212	16251	0.428	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.359	88	13905	2.299	ng/ul#	85
5) Naphthalene	10.801	128	15694	0.391	ng/ul	100
7) 2-Methylnaphthalene	12.407	142	9389	0.370	ng/ul	99
8) 1-Methylnaphthalene	12.627	142	9164	0.349	ng/ul	99
10) Acenaphthylene	14.310	152	13395	0.419	ng/ul	100
11) Acenaphthene	14.643	153	9949	0.391	ng/ul	100
12) Fluorene	15.628	166	10640	0.376	ng/ul	99
14) Pentachlorophenol	16.974	266	2460	0.860	ng/ul	99
15) Phenanthrene	17.375	178	15973	0.402	ng/ul	99
16) Anthracene	17.473	178	14359	0.404	ng/ul	99
19) Fluoranthene	19.394	202	18041	0.394	ng/ul	100
20) Pyrene	19.762	202	18874	0.398	ng/ul	99
21) Benzo(a)anthracene	21.504	228	18334	0.419	ng/ul	99
22) Chrysene	21.560	228	18348	0.415	ng/ul	99
24) Benzo(b)fluoranthene	23.223	252	18867	0.381	ng/ul	94
25) Benzo(k)fluoranthene	23.272	252	20041	0.407	ng/ul	95
26) Benzo(a)pyrene	23.875	252	18298	0.434	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.576	276	23300	0.456	ng/ul	100
28) Dibenzo(a,h)anthracene	26.589	278	18997	0.457	ng/ul	98
29) Benzo(g,h,i)perylene	27.391	276	19134	0.458	ng/ul	98

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

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