

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092523\
 Data File : BM042063.D
 Acq On : 26 Sep 2023 10:15
 Operator : MA/JU
 Sample : PB155715BS
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS715

Quant Time: Sep 27 04:24:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 13:50:58 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	5301	0.400	ng/ul	0.00
4) Naphthalene-d8	10.757	136	12759	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.592	164	6122	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.346	188	12230	0.400	ng/ul	0.00
17) Chrysene-d12	21.518	240	5219	0.400	ng/ul	0.00
23) Perylene-d12	23.924	264	5027	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.368	96	5711	0.791	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.341	152	7882	0.493	ng/ul	-0.01
18) Fluoranthene-d10	19.366	212	11173	0.520	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.402	88	16982	2.234	ng/ul#	85
5) Naphthalene	10.807	128	18294	0.428	ng/ul	100
7) 2-Methylnaphthalene	12.418	142	11441	0.449	ng/ul	99
8) 1-Methylnaphthalene	12.638	142	11061	0.428	ng/ul	100
10) Acenaphthylene	14.319	152	17942	0.437	ng/ul	99
11) Acenaphthene	14.652	153	13199	0.420	ng/ul	100
12) Fluorene	15.642	166	14811	0.418	ng/ul	99
14) Pentachlorophenol	16.978	266	3446	0.533	ng/ul	99
15) Phenanthrene	17.388	178	23008	0.406	ng/ul	99
16) Anthracene	17.481	178	20358	0.415	ng/ul	99
19) Fluoranthene	19.399	202	24293	0.412	ng/ul	99
20) Pyrene	19.766	202	24678	0.426	ng/ul	96
21) Benzo(a)anthracene	21.501	228	15589	0.394	ng/ul	99
22) Chrysene	21.556	228	15853	0.388	ng/ul	99
24) Benzo(b)fluoranthene	23.181	252	14970	0.387	ng/ul	97
25) Benzo(k)fluoranthene	23.231	252	14785	0.393	ng/ul	97
26) Benzo(a)pyrene	23.816	252	13196	0.415	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.418	276	16887	0.387	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.434	278	12382	0.392	ng/ul	99
29) Benzo(g,h,i)perylene	27.192	276	14417	0.392	ng/ul	100

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

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