

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071924\
 Data File : BN032836.D
 Acq On : 19 Jul 2024 12:09
 Operator : MA/JU
 Sample : PB162090BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS090

Quant Time: Jul 19 13:05:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 17 02:35:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.548	152	2733	0.400	ng/ul	0.00
4) Naphthalene-d8	10.321	136	8874	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.193	164	5073	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.965	188	9108	0.400	ng/ul	0.00
17) Chrysene-d12	21.173	240	8151	0.400	ng/ul	0.00
23) Perylene-d12	23.368	264	10047	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	2248	0.685	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.933	152	4099	0.306	ng/ul	0.00
18) Fluoranthene-d10	19.004	212	7786	0.293	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.122	88	6046	1.618	ng/ul	92
5) Naphthalene	10.376	128	7277	0.305	ng/ul	97
7) 2-Methylnaphthalene	12.010	142	4258	0.286	ng/ul	100
8) 1-Methylnaphthalene	12.219	142	4395	0.286	ng/ul	100
10) Acenaphthylene	13.920	152	6233	0.304	ng/ul	97
11) Acenaphthene	14.258	153	5060	0.299	ng/ul	100
12) Fluorene	15.266	166	5018	0.259	ng/ul	99
14) Pentachlorophenol	16.653	266	1081	0.676	ng/ul	100
15) Phenanthrene	17.008	178	7081	0.285	ng/ul	97
16) Anthracene	17.109	178	6153	0.304	ng/ul	97
19) Fluoranthene	19.031	202	8981	0.271	ng/ul	98
20) Pyrene	19.399	202	9538	0.278	ng/ul	99
21) Benzo(a)anthracene	21.159	228	8024	0.309	ng/ul	98
22) Chrysene	21.209	228	10640	0.283	ng/ul	98
24) Benzo(b)fluoranthene	22.708	252	9364	0.264	ng/ul	86
25) Benzo(k)fluoranthene	22.751	252	11184	0.259	ng/ul#	85
26) Benzo(a)pyrene	23.275	252	9509	0.310	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	25.579	276	12529	0.302	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.586	278	9803	0.304	ng/ul#	91
29) Benzo(g,h,i)perylene	26.257	276	10080	0.290	ng/ul	93

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

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