

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092424\
 Data File : BN034214.D
 Acq On : 24 Sep 2024 22:58
 Operator : RC/JU
 Sample : PB163587BS
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS587

Quant Time: Sep 25 00:10:30 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 : Mon Sep 23 19:39:10 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.423	152	5516	0.400	ng/ul	0.00
4) Naphthalene-d8	10.174	136	15297	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.069	164	7930	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.827	188	14053	0.400	ng/ul	0.00
17) Chrysene-d12	21.046	240	9048	0.400	ng/ul	# 0.00
23) Perylene-d12	23.162	264	9285	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.039	96	4547	0.802	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.780	152	7656	0.349	ng/ul	0.00
18) Fluoranthene-d10	18.870	212	11665	0.432	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.069	88	12975	2.062	ng/ul#	78
5) Naphthalene	10.223	128	15152	0.370	ng/ul	100
7) 2-Methylnaphthalene	11.851	142	9315	0.359	ng/ul	100
8) 1-Methylnaphthalene	12.077	142	11171	0.419	ng/ul	100
10) Acenaphthylene	13.782	152	11606	0.322	ng/ul	100
11) Acenaphthene	14.134	153	9941	0.382	ng/ul	98
12) Fluorene	15.133	166	10394	0.358	ng/ul	99
14) Pentachlorophenol	16.485	266	1421	0.321	ng/ul	98
15) Phenanthrene	16.869	178	15192	0.367	ng/ul	100
16) Anthracene	16.962	178	12500	0.335	ng/ul	99
19) Fluoranthene	18.898	202	15217	0.426	ng/ul#	95
20) Pyrene	19.265	202	15682	0.414	ng/ul	98
21) Benzo(a)anthracene	21.031	228	11647	0.327	ng/ul	99
22) Chrysene	21.081	228	14319	0.398	ng/ul	99
24) Benzo(b)fluoranthene	22.539	252	12920	0.358	ng/ul	97
25) Benzo(k)fluoranthene	22.580	252	14751	0.409	ng/ul	97
26) Benzo(a)pyrene	23.074	252	11475	0.397	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.242	276	16032	0.369	ng/ul#	91
28) Dibenzo(a,h)anthracene	25.258	278	12195	0.361	ng/ul	96
29) Benzo(g,h,i)perylene	25.872	276	13257	0.390	ng/ul	96

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

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