

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071624\
 Data File : BN032739.D
 Acq On : 16 Jul 2024 14:03
 Operator : MA/JU
 Sample : PB161904BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS904

Quant Time: Jul 16 14:43:48 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 : Fri Jul 12 01:40:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.553	152	3912	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.332	136	11986	0.400	ng/ul	-0.03
9) Acenaphthene-d10	14.202	164	6951	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.969	188	14167	0.400	ng/ul	-0.02
17) Chrysene-d12	21.185	240	13183	0.400	ng/ul	0.00
23) Perylene-d12	23.383	264	16232	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.101	96	3292	0.701	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.944	152	6471	0.357	ng/ul	-0.03
18) Fluoranthene-d10	19.008	212	14357	0.334	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.131	88	9096	1.700	ng/ul	91
5) Naphthalene	10.382	128	11647	0.362	ng/ul	99
7) 2-Methylnaphthalene	12.015	142	7045	0.351	ng/ul	100
8) 1-Methylnaphthalene	12.230	142	7160	0.345	ng/ul	99
10) Acenaphthylene	13.929	152	10248	0.365	ng/ul	99
11) Acenaphthene	14.267	153	8393	0.361	ng/ul	99
12) Fluorene	15.271	166	8807	0.332	ng/ul	99
14) Pentachlorophenol	16.649	266	1865	0.750	ng/ul	98
15) Phenanthrene	17.012	178	13237	0.343	ng/ul	99
16) Anthracene	17.113	178	11493	0.365	ng/ul	98
19) Fluoranthene	19.041	202	16829	0.314	ng/ul	100
20) Pyrene	19.408	202	17989	0.324	ng/ul	99
21) Benzo(a)anthracene	21.171	228	15652	0.373	ng/ul	99
22) Chrysene	21.220	228	19938	0.328	ng/ul	99
24) Benzo(b)fluoranthene	22.722	252	18452	0.322	ng/ul	96
25) Benzo(k)fluoranthene	22.766	252	20961	0.300	ng/ul	96
26) Benzo(a)pyrene	23.292	252	18416	0.372	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.599	276	24445	0.364	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.606	278	19009	0.365	ng/ul	99
29) Benzo(g,h,i)perylene	26.283	276	19634	0.350	ng/ul	99

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

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