

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102422\
 Data File : BN022291.D
 Acq On : 24 Oct 2022 09:50
 Operator : CG/JU
 Sample : PB148427BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS427

Quant Time: Oct 25 04:06:25 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 22 03:40:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.944	152	3968	0.400	ng/ul	0.00
4) Naphthalene-d8	10.771	136	13637	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.617	164	7331	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.373	188	13000	0.400	ng/ul	0.00
17) Chrysene-d12	21.576	240	6982	0.400	ng/ul	0.00
23) Perylene-d12	24.031	264	5732	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.253	96	3058	0.593	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.366	152	7137	0.346	ng/ul	0.00
18) Fluoranthene-d10	19.410	212	9331	0.398	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.287	88	7620	1.439	ng/ul	88
5) Naphthalene	10.820	128	13071	0.324	ng/ul	100
7) 2-Methylnaphthalene	12.443	142	8251	0.325	ng/ul	99
8) 1-Methylnaphthalene	12.663	142	8859	0.339	ng/ul	100
10) Acenaphthylene	14.340	152	10570	0.321	ng/ul	100
11) Acenaphthene	14.677	153	8856	0.319	ng/ul	99
12) Fluorene	15.667	166	9388	0.291	ng/ul	98
14) Pentachlorophenol	17.027	266	1133	0.367	ng/ul	99
15) Phenanthrene	17.415	178	13664	0.313	ng/ul	99
16) Anthracene	17.508	178	11209	0.305	ng/ul	99
19) Fluoranthene	19.438	202	12177	0.381	ng/ul	99
20) Pyrene	19.805	202	11929	0.371	ng/ul	99
21) Benzo(a)anthracene	21.558	228	8649	0.331	ng/ul	99
22) Chrysene	21.611	228	9309	0.336	ng/ul	98
24) Benzo(b)fluoranthene	23.277	252	9357	0.325	ng/ul	92
25) Benzo(k)fluoranthene	23.327	252	8848	0.316	ng/ul#	93
26) Benzo(a)pyrene	23.923	252	7933	0.346	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.618	276	10303	0.361	ng/ul	100
28) Dibenzo(a,h)anthracene	26.632	278	8066	0.354	ng/ul	97
29) Benzo(g,h,i)perylene	27.413	276	8614	0.371	ng/ul	99

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

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