

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070821\
 Data File : BN015403.D
 Acq On : 08 Jul 2021 10:50
 Operator : CG/JU
 Sample : PB137427BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS427

Quant Time: Jul 08 11:37:54 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 08 11:17:31 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	2834	0.400	ng/ul	0.00
4) Naphthalene-d8	10.485	136	9304	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.340	164	4822	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.086	188	9832	0.400	ng/ul	0.00
17) Chrysene-d12	21.287	240	7950	0.400	ng/ul	0.00
23) Perylene-d12	23.526	264	6869	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	2450	0.767	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.074	152	5205	0.357	ng/ul	0.00
18) Fluoranthene-d10	19.118	212	10534	0.403	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.315	88	6513	1.887	ng/ul	88
5) Naphthalene	10.529	128	10386	0.350	ng/ul	100
7) 2-Methylnaphthalene	12.146	142	6687	0.344	ng/ul	100
8) 1-Methylnaphthalene	12.371	142	6241	0.332	ng/ul	99
10) Acenaphthylene	14.057	152	8150	0.327	ng/ul	99
11) Acenaphthene	14.400	153	6614	0.346	ng/ul	99
12) Fluorene	15.390	166	7342	0.348	ng/ul	99
14) Pentachlorophenol	16.740	266	1310	0.602	ng/ul	99
15) Phenanthrene	17.128	178	12742	0.357	ng/ul	99
16) Anthracene	17.217	178	10647	0.338	ng/ul	100
19) Fluoranthene	19.150	202	13771	0.375	ng/ul	98
20) Pyrene	19.513	202	13852	0.376	ng/ul	98
21) Benzo(a)anthracene	21.270	228	11417	0.388	ng/ul	99
22) Chrysene	21.322	228	13153	0.373	ng/ul	99
24) Benzo(b)fluoranthene	22.857	252	12851	0.400	ng/ul	97
25) Benzo(k)fluoranthene	22.900	252	13135	0.363	ng/ul	98
26) Benzo(a)pyrene	23.430	252	10904	0.405	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	25.774	276	14047	0.432	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.787	278	11208	0.439	ng/ul	98
29) Benzo(g,h,i)perylene	26.458	276	11984	0.397	ng/ul	99

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

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