

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111321\
 Data File : BN017413.D
 Acq On : 14 Nov 2021 05:14
 Operator : CG/JU
 Sample : PB140657BS
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS657

Quant Time: Nov 15 04:39:57 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN111321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 02:31:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.872	152	7109	0.400	ng/ul	0.00
4) Naphthalene-d8	10.667	136	25563	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.502	164	15473	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.247	188	32304	0.400	ng/ul	0.00
17) Chrysene-d12	21.430	240	21681	0.400	ng/ul	0.00
23) Perylene-d12	23.803	264	14395	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.333	96	4751	0.609	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.257	152	12892	0.327	ng/ul	0.00
18) Fluoranthene-d10	19.272	212	28188	0.362	ng/ul	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.366	88	12783	1.271	ng/ul#	50
5) Naphthalene	10.717	128	22110	0.313	ng/ul	97
7) 2-Methylnaphthalene	12.334	142	15460	0.305	ng/ul	99
8) 1-Methylnaphthalene	12.548	142	15063	0.301	ng/ul	100
10) Acenaphthylene	14.220	152	21898	0.295	ng/ul	99
11) Acenaphthene	14.563	153	16976	0.320	ng/ul	98
12) Fluorene	15.553	166	19907	0.316	ng/ul	99
14) Pentachlorophenol	16.893	266	3577	0.506	ng/ul	98
15) Phenanthrene	17.285	178	32157	0.331	ng/ul	99
16) Anthracene	17.378	178	29869	0.306	ng/ul	99
19) Fluoranthene	19.305	202	35603	0.354	ng/ul	100
20) Pyrene	19.663	202	36091	0.351	ng/ul	99
21) Benzo(a)anthracene	21.412	228	26325	0.326	ng/ul	99
22) Chrysene	21.465	228	27901	0.346	ng/ul	99
24) Benzo(b)fluoranthene	23.084	252	21419	0.364	ng/ul	96
25) Benzo(k)fluoranthene	23.131	252	22704	0.375	ng/ul	95
26) Benzo(a)pyrene	23.701	252	17353	0.353	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.259	276	15262	0.368	ng/ul#	87
28) Dibenzo(a,h)anthracene	26.282	278	11489	0.373	ng/ul	97
29) Benzo(g,h,i)perylene	27.010	276	14459	0.381	ng/ul	96

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

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