

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121922\
 Data File : BM038129.D
 Acq On : 20 Dec 2022 07:52
 Operator : CG/JU
 Sample : N6006-17
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Quant Time: Dec 20 23:53:44 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM121522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 17 05:21:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.051	152	5664	0.400	ng/ul	0.00
4) Naphthalene-d8	10.862	136	17930	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.676	164	10356	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.413	188	21820	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.580	240	16352	0.400	ng/ul	0.00
23) Perylene-d12	24.029	264	15229	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.402	96	26664	3.444	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.440	152	7402	0.283	ng/ul	0.00
18) Fluoranthene-d10	19.427	212	16585	0.280	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.911	128	1643	0.030	ng/ul#	85
7) 2-Methylnaphthalene	12.512	142	691	0.021	ng/ul	92
10) Acenaphthylene	14.398	152	1418	0.026	ng/ul#	79
14) Pentachlorophenol	17.063	266	197	0.022	ng/ul	95
15) Phenanthrene	17.451	178	13550	0.184	ng/ul	98
16) Anthracene	17.544	178	2429	0.036	ng/ul#	84
19) Fluoranthene	19.459	202	41498	0.491	ng/ul	100
20) Pyrene	19.822	202	35434m	0.411	ng/ul	
21) Benzo(a)anthracene	21.562	228	20856	0.295	ng/ul	97
22) Chrysene	21.615	228	26379	0.370	ng/ul	96
24) Benzo(b)fluoranthene	23.275	252	36666	0.508	ng/ul	98
25) Benzo(k)fluoranthene	23.322	252	11934m	0.171	ng/ul	
26) Benzo(a)pyrene	23.918	252	23590	0.388	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.582	276	17653	0.231	ng/ul#	85
28) Dibenzo(a,h)anthracene	26.592	278	4781	0.080	ng/ul#	61
29) Benzo(g,h,i)perylene	27.377	276	16685	0.257	ng/ul	97

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

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