

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091523\
 Data File : BN027722.D
 Acq On : 16 Sep 2023 08:26
 Operator : MA/JU
 Sample : PB155388BS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS388

Quant Time: Sep 18 00:49:40 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN091523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 16 01:11:24 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.013	152	4378	0.400	ng/ul	0.00
4) Naphthalene-d8	10.834	136	13817	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.647	164	7928	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.400	188	19485	0.400	ng/ul	0.00
17) Chrysene-d12	21.577	240	17483	0.400	ng/ul	0.00
23) Perylene-d12	24.076	264	16366	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	4235	0.768	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.412	152	8709	0.416	ng/ul	0.00
18) Fluoranthene-d10	19.422	212	23454	0.439	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.431	88	10685	1.862	ng/ul#	85
5) Naphthalene	10.883	128	15140	0.383	ng/ul	100
7) 2-Methylnaphthalene	12.484	142	9491	0.379	ng/ul	99
8) 1-Methylnaphthalene	12.704	142	10403	0.402	ng/ul	100
10) Acenaphthylene	14.379	152	13496	0.390	ng/ul	99
11) Acenaphthene	14.712	153	11645	0.389	ng/ul	99
12) Fluorene	15.698	166	13629	0.387	ng/ul	99
14) Pentachlorophenol	17.033	266	3164	0.694	ng/ul	99
15) Phenanthrene	17.443	178	23878	0.390	ng/ul	99
16) Anthracene	17.535	178	20037	0.391	ng/ul	99
19) Fluoranthene	19.450	202	28079	0.397	ng/ul	99
20) Pyrene	19.817	202	28775	0.395	ng/ul	99
21) Benzo(a)anthracene	21.562	228	26016	0.385	ng/ul	100
22) Chrysene	21.618	228	27263	0.390	ng/ul	100
24) Benzo(b)fluoranthene	23.298	252	27408	0.376	ng/ul	99
25) Benzo(k)fluoranthene	23.351	252	27319	0.386	ng/ul	99
26) Benzo(a)pyrene	23.962	252	22351	0.381	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.706	276	27827	0.370	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.726	278	22402	0.370	ng/ul	99
29) Benzo(g,h,i)perylene	27.525	276	23796	0.373	ng/ul	99

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

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