

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102723\
 Data File : BM042476.D
 Acq On : 27 Oct 2023 14:01
 Operator : MA/JU
 Sample : PB156597TB
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB156597TB

Quant Time: Oct 27 14:34:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 21 02:32:53 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	0.000	152	0	0.000	ng	-8.00
21) Naphthalene-d8	0.000	136	0	0.000	ng	-10.82
39) Acenaphthene-d10	0.000	164	0	0.000	ng	-14.64
64) Phenanthrene-d10	0.000	188	0	0.000	ng	-17.40
76) Chrysene-d12	21.580	240	2063	20.000	ng	# 0.00
86) Perylene-d12	24.038	264	1544	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	1348032	0.000	ng	0.00
7) Phenol-d6	7.151	99	1808159	0.000	ng	0.00
23) Nitrobenzene-d5	9.192	82	1135141	0.000	ng	0.00
42) 2,4,6-Tribromophenol	16.133	330	775968	0.000	ng	0.00
45) 2-Fluorobiphenyl	13.263	172	2502372	0.000	ng	0.00
79) Terphenyl-d14	20.003	244	4963155	42666.563	ng	0.00
Target Compounds						
						Qvalue
77) Benzidine	19.674	184	849	28.452	ng	# 67
78) Pyrene	19.815	202	908	6.299	ng	# 72
80) Butylbenzylphthalate	20.686	149	512	10.562	ng	# 92
81) Benzo(a)anthracene	21.562	228	2170	15.691	ng	# 69
82) 3,3'-Dichlorobenzidine	21.515	252	137	3.318	ng	# 66
83) Chrysene	21.615	228	3699	27.847	ng	# 86
84) Bis(2-ethylhexyl)phtha...	21.433	149	3242	35.817	ng	# 97
85) Di-n-octyl phthalate	22.415	149	251	5.649	ng	# 14
87) Indeno(1,2,3-cd)pyrene	26.632	276	1335	11.831	ng	# 52
88) Benzo(b)fluoranthene	23.280	252	2902	31.225	ng	# 5
89) Benzo(k)fluoranthene	23.321	252	2563	26.563	ng	# 1
90) Benzo(a)pyrene	23.932	252	1377	15.394	ng	# 1
91) Dibenzo(a,h)anthracene	26.632	278	601	6.406	ng	# 1
92) Benzo(g,h,i)perylene	27.432	276	2092	22.962	ng	# 79

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

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